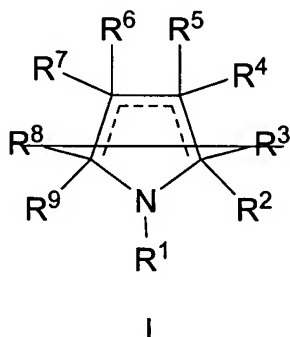
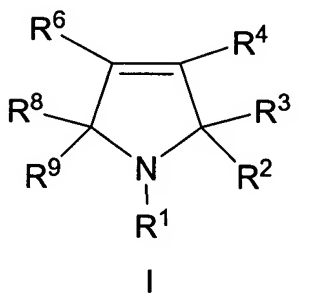


**In the claims:**

1. (Currently amended)



A compound of Formula I:



or a pharmaceutically acceptable salt or stereoisomer thereof, wherein

- a is 0 or 1;
- b is 0 or 1;
- m is 0, 1, or 2;
- n is 0 or 1;
- r is 0 or 1;
- s is 0 or 1;
- u is 2, 3, 4 or 5;

a dashed line represents an optional double bond, provided that one and only one double bond is present in the ring;

R<sup>1</sup> is selected from:

- 1) —(C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)C<sub>1</sub>-C<sub>10</sub>-alkyl,
- 2) —(C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)aryl,
- 3) —(C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)C<sub>2</sub>-C<sub>10</sub>-alkenyl,
- 4) —(C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)C<sub>2</sub>-C<sub>10</sub>-alkynyl,
- 5) —(C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)C<sub>3</sub>-C<sub>8</sub>-cycloalkyl,
- 6) —(C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)heterocyclyl,

- 7)  $(\text{C}_1\text{-C}_6\text{-alkylene})_n(\text{C}=\text{X})\text{NR}^c\text{R}^{c'}$ ,
- 8)  $(\text{C}_1\text{-C}_6\text{-alkylene})_n\text{SO}_2\text{NR}^e\text{R}^{e'}$ ,
- 9)  $(\text{C}_1\text{-C}_6\text{-alkylene})_n\text{SO}_2\text{C}_1\text{-C}_{10}\text{-alkyl}$ ,
- 10)  $(\text{C}_1\text{-C}_6\text{-alkylene})_n\text{SO}_2\text{C}_2\text{-C}_{10}\text{-alkenyl}$ ,
- 11)  $(\text{C}_1\text{-C}_6\text{-alkylene})_n\text{SO}_2\text{C}_2\text{-C}_{10}\text{-alkynyl}$ ,
- 12)  $(\text{C}_1\text{-C}_6\text{-alkylene})_n\text{SO}_2\text{-aryl}$ ,
- 13)  $(\text{C}_1\text{-C}_6\text{-alkylene})_n\text{SO}_2\text{-heterocyclyl}$ ,
- 14)  $(\text{C}_1\text{-C}_6\text{-alkylene})_n\text{SO}_2\text{-C}_3\text{-C}_8\text{-cycloalkyl}$ ,
- 15)  $(\text{C}_1\text{-C}_6\text{-alkylene})_n\text{P}(\text{-O})\text{R}^d\text{R}^{d'}$ ,
- 16)  $\text{aryl}$ ;
- 17)  $\text{heterocyclyl}$ ; and
- 18)  $\text{C}_1\text{-C}_{10}\text{-alkyl}$ ;

said ~~alkyl, aryl, alkenyl, alkynyl, cycloalkyl, alkylene, heteroaryl and heterocyclyl~~ is optionally substituted with one or more substituents selected from R<sup>10</sup>;

R<sup>2</sup> and R<sup>6</sup> are independently selected from:

- 1)  $\text{aryl}$ ,
- 2)  $\text{C}_1\text{-C}_6\text{-aralkyl}$ ,
- 3)  $\text{C}_3\text{-C}_8\text{-cycloalkyl}$ , and
- ~~4)  $\text{heterocyclyl}$ ,~~

said ~~aryl, cycloalkyl, aralkyl and heterocyclyl~~ is optionally substituted with one or more substituents selected from R<sup>10</sup>;

R<sup>3</sup>, R<sup>4</sup>, R<sup>5</sup>, R<sup>7</sup>, R<sup>8</sup>, and R<sup>9</sup> are independently selected from:

- 1)  $\text{H}$ , and
- 2)  $\text{C}_1\text{-C}_{10}\text{ alkyl}$ ,
- 3)  $\text{aryl}$ ,
- 4)  $\text{C}_2\text{-C}_{10}\text{-alkenyl}$ ,
- 5)  $\text{C}_2\text{-C}_{10}\text{-alkynyl}$ ,
- 7)  $\text{C}_1\text{-C}_6\text{-perfluoroalkyl}$ ,
- 8)  $\text{C}_1\text{-C}_6\text{-aralkyl}$ ,

9) ~~—C<sub>3</sub>-C<sub>8</sub> cycloalkyl, and~~

10) ~~—heterocyclyl,~~

~~said alkyl, aryl, alkenyl, alkynyl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>; or~~

~~R<sup>4</sup> and R<sup>5</sup>, or R<sup>8</sup> and R<sup>9</sup>, attached to the same carbon atom are combined to form -(CH<sub>2</sub>)<sub>u</sub> wherein one of the carbon atoms is optionally replaced by a moiety selected from O, S(O)<sub>m</sub>, N(R<sup>a</sup>)C(O), N(R<sup>b</sup>) and N(COR<sup>a</sup>);~~

R<sup>10</sup> is independently selected from:

- 1) (C=O)<sub>a</sub>O<sub>b</sub>C<sub>1</sub>-C<sub>10</sub> alkyl,
- 2) (C=O)<sub>a</sub>O<sub>b</sub>aryl,
- 3) C<sub>2</sub>-C<sub>10</sub> alkenyl,
- 4) C<sub>2</sub>-C<sub>10</sub> alkynyl,
- 5) (C=O)<sub>a</sub>O<sub>b</sub> heterocyclyl,
- 6) CO<sub>2</sub>H,
- 7) halo,
- 8) CN,
- 9) OH,
- 10) O<sub>b</sub>C<sub>1</sub>-C<sub>6</sub> perfluoroalkyl,
- 11) O<sub>a</sub>(C=O)<sub>b</sub>NR<sup>12</sup>R<sup>13</sup>,
- 12) S(O)<sub>m</sub>R<sup>a</sup>,
- 13) S(O)<sub>2</sub>NR<sup>12</sup>R<sup>13</sup>,
- 14) ~~—oxo,~~
- 15) CHO, and
- 16) ~~—(N=O)R<sup>12</sup>R<sup>13</sup>, or~~
- 17) (C=O)<sub>a</sub>O<sub>b</sub>C<sub>3</sub>-C<sub>8</sub> cycloalkyl,

said alkyl, aryl, alkenyl, alkynyl, heterocyclyl, and cycloalkyl optionally substituted with one or more substituents selected from R<sup>11</sup>;

R<sup>11</sup> is selected from:

- 1) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>1</sub>-C<sub>10</sub>)alkyl,

- 2)  $O_r(C_1-C_3)\text{perfluoroalkyl}$ ,
- 3)  $(C_0-C_6)\text{alkylene-S(O)}_mR^a$ ,
- 4) oxo,
- 5) OH,
- 6) halo,
- 7) CN,
- 8)  $(C=O)_rO_s(C_2-C_{10})\text{alkenyl}$ ,
- 9)  $(C=O)_rO_s(C_2-C_{10})\text{alkynyl}$ ,
- 10)  $(C=O)_rO_s(C_3-C_6)\text{cycloalkyl}$ ,
- 11)  $(C=O)_rO_s(C_0-C_6)\text{alkylene-aryl}$ ,
- 12)  $(C=O)_rO_s(C_0-C_6)\text{alkylene-heterocyclyl}$ ,
- 13)  $(C=O)_rO_s(C_0-C_6)\text{alkylene-N(R}^b)_2$ ,
- 14)  $C(O)R^a$ ,
- 15)  $(C_0-C_6)\text{alkylene-CO}_2R^a$ ,
- 16)  $C(O)H$ ,
- 17)  $(C_0-C_6)\text{alkylene-CO}_2H$ ,
- 18)  $C(O)N(R^b)_2$ ,
- 19)  $S(O)_mR^a$ , and
- 20)  $S(O)_2N(R^b)_2$

said alkyl, alkenyl, alkynyl, cycloalkyl, aryl, alkylene and heterocyclyl is optionally substituted with up to three substituents selected from  $R^b$ , OH,  $(C_1-C_6)\text{alkoxy}$ , halogen,  $\text{CO}_2H$ , CN,  $O(C=O)C_1-C_6$  alkyl, oxo, and  $N(R^b)_2$ ;

$R^{12}$  and  $R^{13}$  are independently selected from:

- 1) H,
- 2)  $(C=O)O_bC_1-C_{10}$  alkyl,
- 3)  $(C=O)O_bC_3-C_8$  cycloalkyl,
- 4)  $(C=O)O_b\text{aryl}$ ,
- 5)  $(C=O)O_b\text{heterocyclyl}$ ,
- 6)  $C_1-C_{10}$  alkyl,
- 7) aryl,
- 8)  $C_2-C_{10}$  alkenyl,

- 9) C<sub>2</sub>-C<sub>10</sub> alkynyl,
- 10) heterocyclyl,
- 11) C<sub>3</sub>-C<sub>8</sub> cycloalkyl,
- 12) SO<sub>2</sub>R<sup>a</sup>, and
- 13) (C=O)NR<sup>b</sup><sub>2</sub>,

said alkyl, cycloalkyl, aryl, heterocyclyl, alkenyl, and alkynyl is optionally substituted with one or more substituents selected from R<sup>11</sup>, or

R<sup>12</sup> and R<sup>13</sup> can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 3-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one or more substituents selected from R<sup>11</sup>;

R<sup>14</sup> is independently selected from:

- 1) (C=O)<sub>a</sub>O<sub>b</sub>C<sub>1</sub>-C<sub>10</sub> alkyl,
- 2) (C=O)<sub>a</sub>O<sub>b</sub>aryl,
- 3) C<sub>2</sub>-C<sub>10</sub> alkenyl,
- 4) C<sub>2</sub>-C<sub>10</sub> alkynyl,
- 5) (C=O)<sub>a</sub>O<sub>b</sub> heterocyclyl,
- 6) CO<sub>2</sub>H,
- 7) halo,
- 8) CN,
- 9) OH,
- 10) O<sub>b</sub>C<sub>1</sub>-C<sub>6</sub> perfluoroalkyl,
- 11) O<sub>a</sub>(C=O)<sub>b</sub>NR<sup>12</sup>R<sup>13</sup>,
- 12) S(O)<sub>m</sub>R<sup>a</sup>,
- 13) S(O)<sub>2</sub>NR<sup>12</sup>R<sup>13</sup>,
- 14) oxo,
- 15) CHO,
- 16) (N=O)R<sup>12</sup>R<sup>13</sup>, or
- 17) (C=O)<sub>a</sub>O<sub>b</sub>C<sub>3</sub>-C<sub>8</sub> cycloalkyl,

said alkyl, aryl, alkenyl, alkynyl, heterocyclyl, and cycloalkyl optionally substituted with one or more substituents selected from R<sup>11</sup>;

R<sup>a</sup> is (C<sub>1</sub>-C<sub>6</sub>)alkyl, (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, aryl, or heterocyclyl, optionally substituted with one to three substituents selected from R<sup>14</sup>;

R<sup>b</sup> is H, (C<sub>1</sub>-C<sub>6</sub>)alkyl, aryl, heterocyclyl, (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, (C=O)OC<sub>1</sub>-C<sub>6</sub> alkyl, (C=O)C<sub>1</sub>-C<sub>6</sub> alkyl or S(O)<sub>2</sub>R<sup>a</sup>, optionally substituted with one to three substituents selected from R<sup>14</sup>;

R<sup>c</sup> and R<sup>c'</sup> are independently selected from: H, (C<sub>1</sub>-C<sub>6</sub>)alkyl, aryl, heterocyclyl and (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, optionally substituted with one, two or three substituents selected from R<sup>10</sup>, or R<sup>c</sup> and R<sup>c'</sup> can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 3-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R<sup>11</sup>;

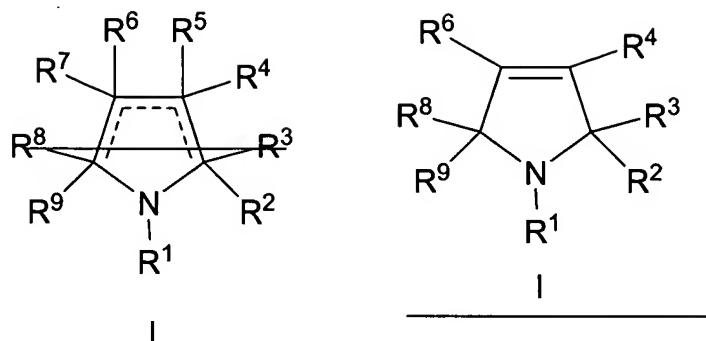
~~R<sup>d</sup> and R<sup>d'</sup> are independently selected from: (C<sub>1</sub>-C<sub>6</sub>)alkyl, (C<sub>1</sub>-C<sub>6</sub>)alkoxy and NR<sup>b</sup><sub>2</sub>, or~~

~~R<sup>d</sup> and R<sup>d'</sup> can be taken together with the phosphorous to which they are attached to form a monocyclic heterocycle with 5-7 members the ring and optionally containing, in addition to the phosphorous, one or two additional heteroatoms selected from NR<sup>e</sup>, O and S, said monocyclic heterocycle optionally substituted with one, two or three substituents selected from R<sup>11</sup>;~~

R<sup>e</sup> is selected from: H and (C<sub>1</sub>-C<sub>6</sub>)alkyl; and

X is selected from O, NR<sup>e</sup> and S.

2. (Currently amended) The compound according to Claim 1 of the Formula I:



or a pharmaceutically acceptable salt or stereoisomer thereof, wherein

- a is 0 or 1;  
b is 0 or 1;  
m is 0, 1, or 2;  
n is 0 or 1;  
r is 0 or 1;  
s is 0 or 1;  
u is 2, 3, 4 or 5;

a dashed line represents an optional double bond, provided that one and only one double bond is present in the ring;

R<sup>1</sup> is selected from:

- 1) —(C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)C<sub>1</sub>-C<sub>10</sub>-alkyl,
- 2) —(C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)aryl,
- 3) —(C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)C<sub>2</sub>-C<sub>10</sub>-alkenyl,
- 4) —(C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)C<sub>2</sub>-C<sub>10</sub>-alkynyl,
- 5) —(C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)C<sub>3</sub>-C<sub>8</sub>-cycloalkyl,
- 6) —(C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)heterocyclyl,
- 7) (C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)NR<sup>c</sup>R<sup>c'</sup>,
- 8) —(C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>SO<sub>2</sub>NR<sup>e</sup>R<sup>e'</sup>,
- 9) —(C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>SO<sub>2</sub>C<sub>1</sub>-C<sub>10</sub>-alkyl,
- 10) —(C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>SO<sub>2</sub>C<sub>2</sub>-C<sub>10</sub>-alkenyl,

- 11) —(C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>SO<sub>2</sub>C<sub>2</sub>-C<sub>10</sub>-alkynyl,
- 12) —(C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>SO<sub>2</sub>-aryl,
- 13) —(C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>SO<sub>2</sub>-heterocyclyl,
- 14) —(C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>SO<sub>2</sub>-C<sub>3</sub>-C<sub>8</sub>-cycloalkyl,
- 15) —(C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>P(=O)R<sup>d</sup>R<sup>d'</sup>,
- 16) —aryl;
- 17) —heterocyclyl; and
- 18) —C<sub>1</sub>-C<sub>10</sub>-alkyl;

said alkyl, aryl, alkenyl, alkynyl, cycloalkyl, alkylene, heteroaryl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;

R<sup>2</sup> and R<sup>6</sup> are independently selected from:

- 1) aryl,
- 2) —C<sub>1</sub>-C<sub>6</sub>-aralkyl,
- 3) —C<sub>3</sub>-C<sub>8</sub>-cycloalkyl, and
- 4) —heterocyclyl,

said aryl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;

R<sup>3</sup>, R<sup>4</sup>, R<sup>5</sup>, R<sup>7</sup>, R<sup>8</sup>, and R<sup>9</sup> are independently selected from:

- 1) H, and
- 2) C<sub>1</sub>-C<sub>10</sub> alkyl,
- 3) —aryl,
- 4) —C<sub>2</sub>-C<sub>10</sub> alkenyl,
- 5) —C<sub>2</sub>-C<sub>10</sub> alkynyl,
- 6) —C<sub>1</sub>-C<sub>6</sub> perfluoroalkyl,
- 7) —C<sub>1</sub>-C<sub>6</sub>-aralkyl,
- 8) —C<sub>3</sub>-C<sub>8</sub>-cycloalkyl, and
- 9) —heterocyclyl,

said alkyl, aryl, alkenyl, alkynyl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>; or



~~R<sup>4</sup> and R<sup>5</sup>, or R<sup>8</sup> and R<sup>9</sup>, attached to the same carbon atom are combined to form  
(CH<sub>2</sub>)<sub>u</sub> wherein one of the carbon atoms is optionally replaced by a moiety selected from O,  
S(O)<sub>m</sub>, N(R<sup>a</sup>)C(O), N(R<sup>b</sup>) and N(COR<sup>a</sup>);~~

R<sup>10</sup> is independently selected from:

- 1) (C=O)<sub>a</sub>O<sub>b</sub>C<sub>1</sub>-C<sub>10</sub> alkyl,
- 2) (C=O)<sub>a</sub>O<sub>b</sub>aryl,
- 3) C<sub>2</sub>-C<sub>10</sub> alkenyl,
- 4) C<sub>2</sub>-C<sub>10</sub> alkynyl,
- 5) (C=O)<sub>a</sub>O<sub>b</sub> heterocyclyl,
- 6) CO<sub>2</sub>H,
- 7) halo,
- 8) CN,
- 9) OH,
- 10) O<sub>b</sub>C<sub>1</sub>-C<sub>6</sub> perfluoroalkyl,
- 11) O<sub>a</sub>(C=O)<sub>b</sub>NR<sup>12</sup>R<sup>13</sup>,
- 12) S(O)<sub>m</sub>R<sup>a</sup>,
- 13) S(O)<sub>2</sub>NR<sup>12</sup>R<sup>13</sup>,
- 14) ~~oxo~~,
- 15) CHO, and
- 16) ~~(N=O)R<sup>12</sup>R<sup>13</sup>, or~~
- 17) (C=O)<sub>a</sub>O<sub>b</sub>C<sub>3</sub>-C<sub>8</sub> cycloalkyl,

said alkyl, aryl, alkenyl, alkynyl, heterocyclyl, and cycloalkyl optionally substituted with one or more substituents selected from R<sup>11</sup>;

R<sup>11</sup> is selected from:

- 1) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>1</sub>-C<sub>10</sub>)alkyl,
- 2) O<sub>r</sub>(C<sub>1</sub>-C<sub>3</sub>)perfluoroalkyl,
- 3) (C<sub>0</sub>-C<sub>6</sub>)alkylene-S(O)<sub>m</sub>R<sup>a</sup>,
- 4) oxo,
- 5) OH,
- 6) halo,

- 7) CN,
- 8)  $(\text{C}=\text{O})_r\text{O}_s(\text{C}_2\text{-C}_{10})\text{alkenyl}$ ,
- 9)  $(\text{C}=\text{O})_r\text{O}_s(\text{C}_2\text{-C}_{10})\text{alkynyl}$ ,
- 10)  $(\text{C}=\text{O})_r\text{O}_s(\text{C}_3\text{-C}_6)\text{cycloalkyl}$ ,
- 11)  $(\text{C}=\text{O})_r\text{O}_s(\text{C}_0\text{-C}_6)\text{alkylene-aryl}$ ,
- 12)  $(\text{C}=\text{O})_r\text{O}_s(\text{C}_0\text{-C}_6)\text{alkylene-heterocyclyl}$ ,
- 13)  $(\text{C}=\text{O})_r\text{O}_s(\text{C}_0\text{-C}_6)\text{alkylene-N(R}^b)_2$ ,
- 14)  $\text{C(O)R}^a$ ,
- 15)  $(\text{C}_0\text{-C}_6)\text{alkylene-CO}_2\text{R}^a$ ,
- 16)  $\text{C(O)H}$ ,
- 17)  $(\text{C}_0\text{-C}_6)\text{alkylene-CO}_2\text{H}$ ,
- 18)  $\text{C(O)N(R}^b)_2$ ,
- 19)  $\text{S(O)}_m\text{R}^a$ , and
- 20)  $\text{S(O)}_2\text{N(R}^b)_2$

said alkyl, alkenyl, alkynyl, cycloalkyl, aryl, alkylene and heterocyclyl is optionally substituted with up to three substituents selected from  $\text{R}^b$ , OH,  $(\text{C}_1\text{-C}_6)\text{alkoxy}$ , halogen,  $\text{CO}_2\text{H}$ , CN,  $\text{O(C=O)C}_1\text{-C}_6$  alkyl, oxo, and  $\text{N(R}^b)_2$ ;

$\text{R}^{12}$  and  $\text{R}^{13}$  are independently selected from:

- 1) H,
- 2)  $(\text{C}=\text{O})\text{O}_b\text{C}_1\text{-C}_{10}$  alkyl,
- 3)  $(\text{C}=\text{O})\text{O}_b\text{C}_3\text{-C}_8$  cycloalkyl,
- 4)  $(\text{C}=\text{O})\text{O}_b\text{aryl}$ ,
- 5)  $(\text{C}=\text{O})\text{O}_b\text{heterocyclyl}$ ,
- 6)  $\text{C}_1\text{-C}_{10}$  alkyl,
- 7) aryl,
- 8)  $\text{C}_2\text{-C}_{10}$  alkenyl,
- 9)  $\text{C}_2\text{-C}_{10}$  alkynyl,
- 10) heterocyclyl,
- 11)  $\text{C}_3\text{-C}_8$  cycloalkyl,
- 12)  $\text{SO}_2\text{R}^a$ , and
- 13)  $(\text{C}=\text{O})\text{NR}^b_2$ ,

said alkyl, cycloalkyl, aryl, heterocyclyl, alkenyl, and alkynyl is optionally substituted with one or more substituents selected from R<sup>11</sup>, or

R<sup>12</sup> and R<sup>13</sup> can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 5-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one or more substituents selected from R<sup>11</sup>;

R<sup>a</sup> is (C<sub>1</sub>-C<sub>6</sub>)alkyl, (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, aryl, or heterocyclyl;

R<sup>b</sup> is H, (C<sub>1</sub>-C<sub>6</sub>)alkyl, aryl, heterocyclyl, (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, (C=O)OC<sub>1</sub>-C<sub>6</sub> alkyl, (C=O)C<sub>1</sub>-C<sub>6</sub> alkyl or S(O)<sub>2</sub>R<sup>a</sup>;

R<sup>c</sup> and R<sup>c'</sup> are independently selected from: H, (C<sub>1</sub>-C<sub>6</sub>)alkyl, aryl, heterocyclyl and (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, optionally substituted with one, two or three substituents selected from R<sup>10</sup>, or

R<sup>c</sup> and R<sup>c'</sup> can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 5-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R<sup>11</sup>; and

~~R<sup>d</sup> and R<sup>d'</sup> are independently selected from: (C<sub>1</sub>-C<sub>6</sub>)alkyl, (C<sub>1</sub>-C<sub>6</sub>)alkoxy and NR<sup>b</sup><sub>2</sub>, or~~

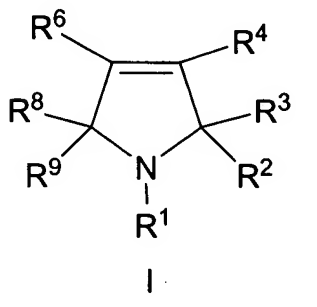
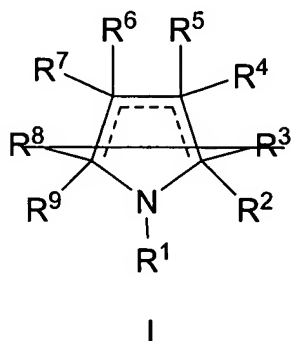
~~R<sup>d</sup> and R<sup>d'</sup> can be taken together with the phosphorous to which they are attached to form a monocyclic heterocycle with 5-7 members the ring and optionally containing, in addition to the phosphorous, one or two additional heteroatoms selected from NR<sup>e</sup>, O and S, said monocyclic heterocycle optionally substituted with one, two or three substituents selected from R<sup>11</sup>;~~

R<sup>e</sup> is selected from: H and (C<sub>1</sub>-C<sub>6</sub>)alkyl; and

X is selected from O, NR<sup>e</sup> and S.

3. (Currently amended)

The compound according to Claim 2 of Formula I:



or a pharmaceutically acceptable salt or stereoisomer thereof, wherein

- a is 0 or 1;
- b is 0 or 1;
- m is 0, 1, or 2;
- n is 0 or 1;
- r is 0 or 1;
- s is 0 or 1;
- u is 2, 3, 4 or 5;

~~a dashed line represents an optional double bond, provided that one and only one double bond is present in the ring;~~

R<sup>1</sup> is selected from:

- 1) —(C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)C<sub>1</sub>-C<sub>10</sub>-alkyl;
- 2) —(C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)aryl;
- 3) —(C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)C<sub>2</sub>-C<sub>10</sub>-alkenyl;
- 4) —(C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)C<sub>2</sub>-C<sub>10</sub>-alkynyl;
- 5) —(C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)C<sub>3</sub>-C<sub>8</sub>-cycloalkyl;
- 6) —(C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=X)heterocyclyl;

- 7)  $(\text{C}_1\text{-C}_6\text{-alkylene})_n(\text{C}=\text{X})\text{NR}^c\text{R}^c$ ,
- 8)  $(\text{C}_1\text{-C}_6\text{-alkylene})_n\text{SO}_2\text{NR}^e\text{R}^e$ ,
- 9)  $(\text{C}_1\text{-C}_6\text{-alkylene})_n\text{SO}_2\text{C}_1\text{-C}_{10}\text{-alkyl}$ ,
- 10)  $(\text{C}_1\text{-C}_6\text{-alkylene})_n\text{SO}_2\text{C}_2\text{-C}_{10}\text{-alkenyl}$ ,
- 11)  $(\text{C}_1\text{-C}_6\text{-alkylene})_n\text{SO}_2\text{C}_2\text{-C}_{10}\text{-alkynyl}$ ,
- 12)  $(\text{C}_1\text{-C}_6\text{-alkylene})_n\text{SO}_2\text{-aryl}$ ,
- 13)  $(\text{C}_1\text{-C}_6\text{-alkylene})_n\text{SO}_2\text{-heterocyclyl}$ ,
- 14)  $(\text{C}_1\text{-C}_6\text{-alkylene})_n\text{SO}_2\text{-C}_3\text{-C}_8\text{-cycloalkyl}$ ,
- 15)  $(\text{C}_1\text{-C}_6\text{-alkylene})_n\text{P}(=\text{O})\text{R}^d\text{R}^d$ ,
- 16)  $\text{aryl}$ ;
- 17)  $\text{heterocyclyl}$ ; and
- 18)  $\text{C}_1\text{-C}_{10}\text{-alkyl}$ ;

said  $\text{alkyl}$ ,  $\text{aryl}$ ,  $\text{alkenyl}$ ,  $\text{alkynyl}$ ,  $\text{cycloalkyl}$ ,  $\text{alkylene}$ ,  $\text{heteroaryl}$  and  $\text{heterocyclyl}$  is optionally substituted with one or more substituents selected from  $\text{R}^{10}$ ;

$\text{R}^2$  and  $\text{R}^6$  are independently selected from:

- 1)  $\text{aryl}$ ,
- 2)  $\text{C}_1\text{-C}_6\text{-aralkyl}$ ,
- 3)  $\text{C}_3\text{-C}_8\text{-cycloalkyl}$ , and
- 4)  $\text{heterocyclyl}$ ,

said  $\text{aryl}$ ,  $\text{cycloalkyl}$ ,  $\text{aralkyl}$  and  $\text{heterocyclyl}$  is optionally substituted with one or more substituents selected from  $\text{R}^{10}$ ;

$\text{R}^3$ ,  $\text{R}^4$ ,  $\text{R}^5$ ,  $\text{R}^7$ ,  $\text{R}^8$ , and  $\text{R}^9$  are independently selected from:

- 1)  $\text{H}$ , and
- 2)  $\text{C}_1\text{-C}_{10}\text{-alkyl}$ ,
- 3)  $\text{aryl}$ ,
- 4)  $\text{C}_2\text{-C}_{10}\text{-alkenyl}$ ,
- 5)  $\text{C}_2\text{-C}_{10}\text{-alkynyl}$ ,
- 6)  $\text{C}_1\text{-C}_6\text{-perfluoroalkyl}$ ,
- 7)  $\text{C}_1\text{-C}_6\text{-aralkyl}$ ,
- 8)  $\text{C}_3\text{-C}_8\text{-cycloalkyl}$ , and

~~—9)—heterocyclyl,~~

~~said alkyl, aryl, alkenyl, alkynyl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>; or~~

~~R<sup>4</sup> and R<sup>5</sup>, or R<sup>8</sup> and R<sup>9</sup>, attached to the same carbon atom are combined to form (CH<sub>2</sub>)<sub>u</sub> wherein one of the carbon atoms is optionally replaced by a moiety selected from O, S(O)<sub>m</sub>, N(R<sup>a</sup>)C(O), N(R<sup>b</sup>) and N(COR<sup>a</sup>);~~

R<sup>10</sup> is independently selected from:

- 1) (C=O)<sub>a</sub>O<sub>b</sub>C<sub>1</sub>-C<sub>10</sub> alkyl,
- 2) (C=O)<sub>a</sub>O<sub>b</sub>aryl,
- 3) C<sub>2</sub>-C<sub>10</sub> alkenyl,
- 4) C<sub>2</sub>-C<sub>10</sub> alkynyl,
- 5) (C=O)<sub>a</sub>O<sub>b</sub> heterocyclyl,
- 6) CO<sub>2</sub>H,
- 7) halo,
- 8) CN,
- 9) OH,
- 10) O<sub>b</sub>C<sub>1</sub>-C<sub>6</sub> perfluoroalkyl,
- 11) O<sub>a</sub>(C=O)<sub>b</sub>NR<sup>12</sup>R<sup>13</sup>,
- 12) S(O)<sub>m</sub>R<sup>a</sup>,
- 13) S(O)<sub>2</sub>NR<sup>12</sup>R<sup>13</sup>,
- ~~14) —oxo,~~
- 15) CHO, and
- ~~16) —(N=O)R<sup>12</sup>R<sup>13</sup>, or~~
- 17) (C=O)<sub>a</sub>O<sub>b</sub>C<sub>3</sub>-C<sub>8</sub> cycloalkyl,

said alkyl, aryl, alkenyl, alkynyl, heterocyclyl, and cycloalkyl optionally substituted with one or more substituents selected from R<sup>11</sup>;

R<sup>11</sup> is selected from:

- 1) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>1</sub>-C<sub>10</sub>)alkyl,
- 2) O<sub>r</sub>(C<sub>1</sub>-C<sub>3</sub>)perfluoroalkyl,
- 3) (C<sub>0</sub>-C<sub>6</sub>)alkylene-S(O)<sub>m</sub>R<sup>a</sup>,

- 4) oxo,
- 5) OH,
- 6) halo,
- 7) CN,
- 8)  $(\text{C}=\text{O})_r\text{O}_s(\text{C}_2\text{-C}_{10})\text{alkenyl}$ ,
- 9)  $(\text{C}=\text{O})_r\text{O}_s(\text{C}_2\text{-C}_{10})\text{alkynyl}$ ,
- 10)  $(\text{C}=\text{O})_r\text{O}_s(\text{C}_3\text{-C}_6)\text{cycloalkyl}$ ,
- 11)  $(\text{C}=\text{O})_r\text{O}_s(\text{C}_0\text{-C}_6)\text{alkylene-aryl}$ ,
- 12)  $(\text{C}=\text{O})_r\text{O}_s(\text{C}_0\text{-C}_6)\text{alkylene-heterocyclyl}$ ,
- 13)  $(\text{C}=\text{O})_r\text{O}_s(\text{C}_0\text{-C}_6)\text{alkylene-N(R}^b)_2$ ,
- 14)  $\text{C(O)R}^a$ ,
- 15)  $(\text{C}_0\text{-C}_6)\text{alkylene-CO}_2\text{R}^a$ ,
- 16)  $\text{C(O)H}$ ,
- 17)  $(\text{C}_0\text{-C}_6)\text{alkylene-CO}_2\text{H}$ ,
- 18)  $\text{C(O)N(R}^b)_2$ ,
- 19)  $\text{S(O)}_m\text{R}^a$ , and
- 20)  $\text{S(O)}_2\text{N(R}^b)_2$ ,

said alkyl, alkenyl, alkynyl, cycloalkyl, aryl, alkylene and heterocyclyl is optionally substituted with up to three substituents selected from  $\text{R}^b$ , OH,  $(\text{C}_1\text{-C}_6)\text{alkoxy}$ , halogen,  $\text{CO}_2\text{H}$ , CN,  $\text{O(C=O)C}_1\text{-C}_6$  alkyl, oxo, and  $\text{N(R}^b)_2$ ;

$\text{R}^{12}$  and  $\text{R}^{13}$  are independently selected from:

- 1) H,
- 2)  $(\text{C}=\text{O})\text{O}_b\text{C}_1\text{-C}_{10}$  alkyl,
- 3)  $(\text{C}=\text{O})\text{O}_b\text{C}_3\text{-C}_8$  cycloalkyl,
- 4)  $(\text{C}=\text{O})\text{O}_b\text{aryl}$ ,
- 5)  $(\text{C}=\text{O})\text{O}_b\text{heterocyclyl}$ ,
- 6)  $\text{C}_1\text{-C}_{10}$  alkyl,
- 7) aryl,
- 8)  $\text{C}_2\text{-C}_{10}$  alkenyl,
- 9)  $\text{C}_2\text{-C}_{10}$  alkynyl,
- 10) heterocyclyl,

- 11) C<sub>3</sub>-C<sub>8</sub> cycloalkyl,
- 12) SO<sub>2</sub>R<sup>a</sup>, and
- 13) (C=O)NR<sup>b</sup><sub>2</sub>,

said alkyl, cycloalkyl, aryl, heterocyclyl, alkenyl, and alkynyl is optionally substituted with one or more substituents selected from R<sup>11</sup>, or

R<sup>12</sup> and R<sup>13</sup> can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 5-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one or more substituents selected from R<sup>11</sup>;

R<sup>a</sup> is (C<sub>1</sub>-C<sub>6</sub>)alkyl, (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, aryl, or heterocyclyl;

R<sup>b</sup> is H, (C<sub>1</sub>-C<sub>6</sub>)alkyl, aryl, heterocyclyl, (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, (C=O)OC<sub>1</sub>-C<sub>6</sub> alkyl, (C=O)C<sub>1</sub>-C<sub>6</sub> alkyl or S(O)<sub>2</sub>R<sup>a</sup>;

R<sup>c</sup> and R<sup>c'</sup> are independently selected from: H, (C<sub>1</sub>-C<sub>6</sub>)alkyl, aryl, heterocyclyl and (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl or

R<sup>c</sup> and R<sup>c'</sup> can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 5-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R<sup>11</sup>;

~~R<sup>d</sup> and R<sup>d'</sup> are independently selected from: (C<sub>1</sub>-C<sub>6</sub>)alkyl, (C<sub>1</sub>-C<sub>6</sub>)alkoxy and NR<sup>b</sup><sub>2</sub>, or~~

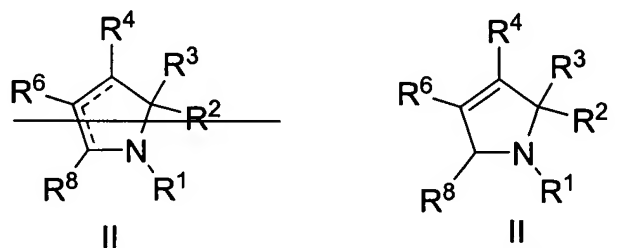
~~R<sup>d</sup> and R<sup>d'</sup> can be taken together with the phosphorous to which they are attached to form a monocyclic heterocycle with 5-7 members the ring and optionally containing, in addition to the phosphorous, one or two additional heteroatoms selected from NR<sup>e</sup>, O and S, said monocyclic heterocycle optionally substituted with one, two or three substituents selected from R<sup>11</sup>;~~



$R^e$  is selected from: H and (C<sub>1</sub>-C<sub>6</sub>)alkyl; and

X is selected from O, NR<sup>e</sup> and S.

4. (Currently amended) The compound according to Claim 2 of the Formula II,



or a pharmaceutically acceptable salt or stereoisomer thereof, wherein  
wherein:

a is 0 or 1;  
b is 0 or 1;  
m is 0, 1, or 2;  
r is 0 or 1;  
s is 0 or 1;

a dashed line represents an optional double bond, provided that one and only one double bond is present in the ring;

R<sup>1</sup> is selected from:

- 1) —(C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=O)C<sub>1</sub>-C<sub>10</sub>-alkyl,
- 2) —(C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=O)aryl,
- 3) —(C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=O)C<sub>2</sub>-C<sub>10</sub>-alkenyl,
- 4) —(C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=O)C<sub>2</sub>-C<sub>10</sub>-alkynyl,
- 5) —(C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=O)C<sub>3</sub>-C<sub>8</sub>-cycloalkyl,
- 6) —(C<sub>1</sub>-C<sub>6</sub>-alkylene)<sub>n</sub>(C=O)heterocyclyl,

- 7)  $(\text{C}_1\text{-C}_6\text{-alkylene})_n(\text{C}=\text{O})\text{NR}^c\text{R}^c$ ,
- 8)  $(\text{C}_1\text{-C}_6\text{-alkylene})_n\text{SO}_2\text{NR}^e\text{R}^{e2}$ ,
- 9)  $(\text{C}_1\text{-C}_6\text{-alkylene})_n\text{SO}_2\text{C}_1\text{-C}_{10}\text{-alkyl}$ ,
- 10)  $(\text{C}_1\text{-C}_6\text{-alkylene})_n\text{SO}_2\text{-aryl}$ ,
- 11)  $(\text{C}_1\text{-C}_6\text{-alkylene})_n\text{SO}_2\text{-heterocyclyl}$ ,
- 12)  $(\text{C}_1\text{-C}_6\text{-alkenyl})_n\text{SO}_2\text{-C}_3\text{-C}_8\text{-cycloalkyl}$ ,
- 13)  $(\text{C}_1\text{-C}_6\text{-alkylene})_n\text{P}(=\text{O})\text{R}^d\text{R}^{d2}$ ,
- 14)  $\text{-aryl}$ ,
- 15)  $\text{-heterocyclyl}$ , and
- 16)  $\text{-C}_1\text{-C}_{10}\text{-alkyl}$ ;

said ~~alkyl, aryl, alkenyl, alkynyl, cycloalkyl, alkylene, heteroaryl and heterocyclyl~~ is optionally substituted with one or more substituents selected from R<sup>10</sup>;

R<sup>2</sup> and R<sup>6</sup> are independently selected from:

- 1)  $\text{aryl}$ ,
- 2)  $\text{-C}_1\text{-C}_6\text{-aralkyl}$ ,
- 3)  $\text{-C}_3\text{-C}_8\text{-cycloalkyl}$ , and
- ~~4)  $\text{-heterocyclyl}$ ,~~

said ~~aryl, cycloalkyl, aralkyl and heterocyclyl~~ is optionally substituted with one or more substituents selected from R<sup>10</sup>;

R<sup>3</sup>, R<sup>4</sup> and R<sup>8</sup> are independently selected from:

- 1)  $\text{H}$ , and
- 2)  $\text{C}_1\text{-C}_{10}\text{ alkyl}$ ,
- 3)  $\text{-aryl}$ ,
- 4)  $\text{-C}_2\text{-C}_{10}\text{-alkenyl}$ ,
- 5)  $\text{-C}_2\text{-C}_{10}\text{-alkynyl}$ ,
- 6)  $\text{-C}_1\text{-C}_6\text{-perfluoroalkyl}$ ,
- 7)  $\text{-C}_1\text{-C}_6\text{-aralkyl}$ ,
- 8)  $\text{-C}_3\text{-C}_8\text{-cycloalkyl}$ , and
- 9)  $\text{-heterocyclyl}$ ,

~~said alkyl, aryl, alkenyl, alkynyl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;~~

R<sup>10</sup> is independently selected from:

- 1) (C=O)<sub>a</sub>O<sub>b</sub>C<sub>1</sub>-C<sub>10</sub> alkyl,
- 2) (C=O)<sub>a</sub>O<sub>b</sub>aryl,
- 3) C<sub>2</sub>-C<sub>10</sub> alkenyl,
- 4) C<sub>2</sub>-C<sub>10</sub> alkynyl,
- 5) (C=O)<sub>a</sub>O<sub>b</sub> heterocyclyl,
- 6) CO<sub>2</sub>H,
- 7) halo,
- 8) CN,
- 9) OH,
- 10) O<sub>b</sub>C<sub>1</sub>-C<sub>6</sub> perfluoroalkyl,
- 11) O<sub>a</sub>(C=O)<sub>b</sub>NR<sup>12</sup>R<sup>13</sup>,
- 12) S(O)<sub>m</sub>R<sup>a</sup>,
- 13) S(O)<sub>2</sub>NR<sup>12</sup>R<sup>13</sup>,
- ~~14) —oxo,~~
- 15) CHO, and
- ~~16) (N=O)R<sup>12</sup>R<sup>13</sup>, or~~
- 17) (C=O)<sub>a</sub>O<sub>b</sub>C<sub>3</sub>-C<sub>8</sub> cycloalkyl,

said alkyl, aryl, alkenyl, alkynyl, heterocyclyl, and cycloalkyl optionally substituted with one, two or three substituents selected from R<sup>11</sup>;

R<sup>11</sup> is selected from:

- 1) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>1</sub>-C<sub>10</sub>)alkyl,
- 2) O<sub>r</sub>(C<sub>1</sub>-C<sub>3</sub>)perfluoroalkyl,
- 3) oxo,
- 4) OH,
- 5) halo,
- 6) CN,
- 7) (C<sub>2</sub>-C<sub>10</sub>)alkenyl,

- 8) (C<sub>2</sub>-C<sub>10</sub>)alkynyl,
- 9) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>3</sub>-C<sub>6</sub>)cycloalkyl,
- 10) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>0</sub>-C<sub>6</sub>)alkylene-aryl,
- 11) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>0</sub>-C<sub>6</sub>)alkylene-heterocyclyl,
- 12) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>0</sub>-C<sub>6</sub>)alkylene-N(R<sup>b</sup>)<sub>2</sub>,
- 13) C(O)R<sup>a</sup>,
- 14) (C<sub>0</sub>-C<sub>6</sub>)alkylene-CO<sub>2</sub>R<sup>a</sup>,
- 15) C(O)H,
- 16) (C<sub>0</sub>-C<sub>6</sub>)alkylene-CO<sub>2</sub>H,
- 17) C(O)N(R<sup>b</sup>)<sub>2</sub>,
- 18) S(O)<sub>m</sub>R<sup>a</sup>, and
- 19) S(O)<sub>2</sub>N(R<sup>b</sup>)<sub>2</sub>;

said alkyl, alkenyl, alkynyl, cycloalkyl, aryl, alkylene and heterocyclyl is optionally substituted with up to three substituents selected from R<sup>b</sup>, OH, (C<sub>1</sub>-C<sub>6</sub>)alkoxy, halogen, CO<sub>2</sub>H, CN, O(C=O)C<sub>1</sub>-C<sub>6</sub> alkyl, oxo, and N(R<sup>b</sup>)<sub>2</sub>;

R<sup>12</sup> and R<sup>13</sup> are independently selected from:

- 1) H,
- 2) (C=O)O<sub>b</sub>C<sub>1</sub>-C<sub>10</sub> alkyl,
- 3) (C=O)O<sub>b</sub>C<sub>3</sub>-C<sub>8</sub> cycloalkyl,
- 4) (C=O)O<sub>b</sub>aryl,
- 5) (C=O)O<sub>b</sub>heterocyclyl,
- 6) C<sub>1</sub>-C<sub>10</sub> alkyl,
- 7) aryl,
- 8) C<sub>2</sub>-C<sub>10</sub> alkenyl,
- 9) C<sub>2</sub>-C<sub>10</sub> alkynyl,
- 10) heterocyclyl,
- 11) C<sub>3</sub>-C<sub>8</sub> cycloalkyl,
- 12) SO<sub>2</sub>R<sup>a</sup>, and
- 13) (C=O)NR<sup>b</sup><sub>2</sub>,

said alkyl, cycloalkyl, aryl, heterocyclyl, alkenyl, and alkynyl is optionally substituted with one, two or three substituents selected from R<sup>11</sup>, or

R<sup>12</sup> and R<sup>13</sup> can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 5-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R<sup>11</sup>;

R<sup>a</sup> is (C<sub>1</sub>-C<sub>6</sub>)alkyl, (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, aryl, or heterocyclyl;

R<sup>b</sup> is H, (C<sub>1</sub>-C<sub>6</sub>)alkyl, aryl, heterocyclyl, (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, (C=O)OC<sub>1</sub>-C<sub>6</sub> alkyl, (C=O)C<sub>1</sub>-C<sub>6</sub> alkyl or S(O)<sub>2</sub>R<sup>a</sup>;

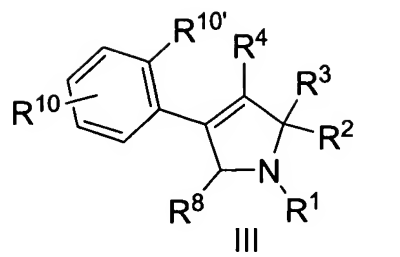
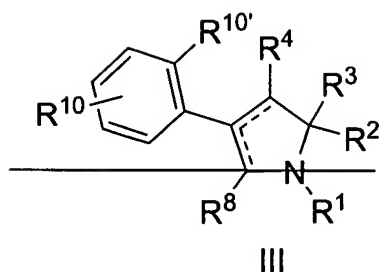
R<sup>c</sup> and R<sup>c'</sup> are independently selected from: H, (C<sub>1</sub>-C<sub>6</sub>)alkyl, aryl, heterocyclyl and (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl; or R<sup>c</sup> and R<sup>c'</sup> can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 5-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R<sup>11</sup>;

~~R<sup>d</sup> and R<sup>d'</sup> are independently selected from: (C<sub>1</sub>-C<sub>6</sub>)alkyl, (C<sub>1</sub>-C<sub>6</sub>)alkoxy and NR<sup>b</sup><sub>2</sub>, or~~

~~R<sup>d</sup> and R<sup>d'</sup> can be taken together with the phosphorous to which they are attached to form a monocyclic heterocycle with 5-7 members the ring and optionally containing, in addition to the phosphorous, one or two additional heteroatoms selected from NR<sup>e</sup>, O and S, said monocyclic heterocycle optionally substituted with one, two or three substituents selected from R<sup>11</sup>; and~~

~~R<sup>e</sup> is selected from: H and (C<sub>1</sub>-C<sub>6</sub>)alkyl.~~

5. (Currently amended)      A compound of the Formula III,



or a pharmaceutically acceptable salt or stereoisomer thereof, wherein

- a is 0 or 1;
- b is 0 or 1;
- m is 0, 1, or 2;
- r is 0 or 1;
- s is 0 or 1;

~~a dashed line represents an optional double bond, provided that one and only one double bond is present in the ring;~~

R<sup>1</sup> is selected from:

- 1) —(C=O)C<sub>1</sub>-C<sub>10</sub>-alkyl,
- 2) —(C=O)aryl,
- 3) —(C=O)C<sub>2</sub>-C<sub>10</sub>-alkenyl,
- 4) —(C=O)C<sub>2</sub>-C<sub>10</sub>-alkynyl,
- 5) —(C=O)C<sub>3</sub>-C<sub>8</sub>-cycloalkyl,
- 6) —(C=O)heterocyclyl,
- 7) (C=O)NR<sup>c</sup>R<sup>c'</sup>,
- 8) —SO<sub>2</sub>NR<sup>e</sup>R<sup>e'</sup>,
- 9) —SO<sub>2</sub>C<sub>1</sub>-C<sub>10</sub>-alkyl,
- 10) —SO<sub>2</sub>-aryl,
- 11) —SO<sub>2</sub>-heterocyclyl,
- 12) —SO<sub>2</sub>-C<sub>3</sub>-C<sub>8</sub>-cycloalkyl, and
- 13) —P(=O)R<sup>d</sup>R<sup>d'</sup>,

~~said alkyl, aryl, alkenyl, alkynyl, cycloalkyl, heteroaryl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;~~

R<sup>2</sup> is selected from:

- 1) aryl,
- 2) ~~C<sub>1</sub>-C<sub>6</sub> aralkyl,~~
- 3) ~~C<sub>3</sub>-C<sub>8</sub> cycloalkyl, and~~
- 4) ~~heterocyclyl,~~

~~said aryl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;~~

R<sup>3</sup>, R<sup>4</sup> and R<sup>8</sup> are independently selected from:

- 1) H, and
- 2) C<sub>1</sub>-C<sub>10</sub> alkyl,
- 3) ~~aryl,~~
- 4) ~~C<sub>2</sub>-C<sub>10</sub> alkenyl,~~
- 5) ~~C<sub>2</sub>-C<sub>10</sub> alkynyl,~~
- 6) ~~C<sub>1</sub>-C<sub>6</sub> perfluoroalkyl,~~
- 7) ~~C<sub>1</sub>-C<sub>6</sub> aralkyl,~~
- 8) ~~C<sub>3</sub>-C<sub>8</sub> cycloalkyl, and~~
- 9) ~~heterocyclyl,~~

~~said alkyl, aryl, alkenyl, alkynyl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;~~

R<sup>10</sup> is independently selected from:

- 1) (C=O)<sub>a</sub>O<sub>b</sub>C<sub>1</sub>-C<sub>10</sub> alkyl,
- 2) (C=O)<sub>a</sub>O<sub>b</sub>aryl,
- 3) C<sub>2</sub>-C<sub>10</sub> alkenyl,
- 4) C<sub>2</sub>-C<sub>10</sub> alkynyl,
- 5) (C=O)<sub>a</sub>O<sub>b</sub> heterocyclyl,
- 6) CO<sub>2</sub>H,
- 7) halo,
- 8) CN,

- 9) OH,
- 10)  $O_bC_1-C_6$  perfluoroalkyl,
- 11)  $O_a(C=O)_bNR^{12}R^{13}$ ,
- 12)  $S(O)_mR^a$ ,
- 13)  $S(O)_2NR^{12}R^{13}$ ,
- ~~14) —oxo,~~
- 15) CHO, and
- ~~16) —(N=O)R<sup>12</sup>R<sup>13</sup>, or~~
- 17)  $(C=O)_aO_bC_3-C_8$  cycloalkyl,

said alkyl, aryl, alkenyl, alkynyl, heterocyclyl, and cycloalkyl optionally substituted with one, two or three substituents selected from R<sup>11</sup>;

R<sup>10'</sup> is halogen;

R<sup>11</sup> is selected from:

- 1)  $(C=O)_rO_s(C_1-C_{10})$ alkyl,
- 2)  $O_r(C_1-C_3)$ perfluoroalkyl,
- 3) oxo,
- 4) OH,
- 5) halo,
- 6) CN,
- 7)  $(C_2-C_{10})$ alkenyl,
- 8)  $(C_2-C_{10})$ alkynyl,
- 9)  $(C=O)_rO_s(C_3-C_6)$ cycloalkyl,
- 10)  $(C=O)_rO_s(C_0-C_6)$ alkylene-aryl,
- 11)  $(C=O)_rO_s(C_0-C_6)$ alkylene-heterocyclyl,
- 12)  $(C=O)_rO_s(C_0-C_6)$ alkylene-N(R<sup>b</sup>)<sub>2</sub>,
- 13) C(O)R<sup>a</sup>,
- 14)  $(C_0-C_6)$ alkylene-CO<sub>2</sub>R<sup>a</sup>,
- 15) C(O)H,
- 16)  $(C_0-C_6)$ alkylene-CO<sub>2</sub>H, and
- 17) C(O)N(R<sup>b</sup>)<sub>2</sub>,



- 18)  $S(O)_mR^a$ , and
- 19)  $S(O)_2N(R^b)_2$ ;

said alkyl, alkenyl, alkynyl, cycloalkyl, aryl, and heterocyclyl is optionally substituted with up to three substituents selected from  $R^b$ , OH,  $(C_1-C_6)$ alkoxy, halogen,  $CO_2H$ , CN,  $O(C=O)C_1-C_6$  alkyl, oxo, and  $N(R^b)_2$ ;

$R^{12}$  and  $R^{13}$  are independently selected from:

- 1) H,
- 2)  $(C=O)O_bC_1-C_{10}$  alkyl,
- 3)  $(C=O)O_bC_3-C_8$  cycloalkyl,
- 4)  $(C=O)O_b$ aryl,
- 5)  $(C=O)O_b$ heterocyclyl,
- 6)  $C_1-C_{10}$  alkyl,
- 7) aryl,
- 8)  $C_2-C_{10}$  alkenyl,
- 9)  $C_2-C_{10}$  alkynyl,
- 10) heterocyclyl,
- 11)  $C_3-C_8$  cycloalkyl,
- 12)  $SO_2R^a$ , and
- 13)  $(C=O)NR^b_2$ ,

said alkyl, cycloalkyl, aryl, heterocyclyl, alkenyl, and alkynyl is optionally substituted with one, two or three substituents selected from  $R^{11}$ , or

$R^{12}$  and  $R^{13}$  can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 5-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from  $R^{11}$ ;

$R^a$  is  $(C_1-C_6)$ alkyl,  $(C_3-C_6)$ cycloalkyl, aryl, or heterocyclyl;

$R^b$  is H, (C<sub>1</sub>-C<sub>6</sub>)alkyl, aryl, heterocyclyl, (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, (C=O)OC<sub>1</sub>-C<sub>6</sub> alkyl, (C=O)C<sub>1</sub>-C<sub>6</sub> alkyl or S(O)<sub>2</sub>R<sub>a</sub>;

$R^c$  and  $R^{c'}$  are independently selected from: H, (C<sub>1</sub>-C<sub>6</sub>)alkyl, aryl, heterocyclyl and (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl; or

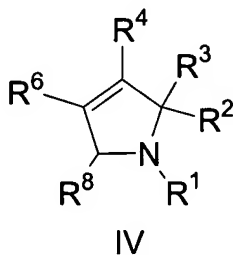
$R^c$  and  $R^{c'}$  can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 5-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from  $R^{11}$ ;

~~$R^d$  and  $R^{d'}$  are independently selected from: (C<sub>1</sub>-C<sub>6</sub>)alkyl, (C<sub>1</sub>-C<sub>6</sub>)alkoxy and  $NR^b$ , or~~

~~$R^d$  and  $R^{d'}$  can be taken together with the phosphorous to which they are attached to form a monocyclic heterocycle with 5-7 members the ring and optionally containing, in addition to the phosphorous, one or two additional heteroatoms selected from  $NR^e$ , O and S, said monocyclic heterocycle optionally substituted with one, two or three substituents selected from  $R^{11}$ ; and~~

~~$R^e$  is selected from: H and (C<sub>1</sub>-C<sub>6</sub>)alkyl.~~

6. (Currently amended) The compound according to Claim 4 of the Formula IV,



or a pharmaceutically acceptable salt or stereoisomer thereof, wherein

a is 0 or 1;  
b is 0 or 1;  
m is 0, 1, or 2;  
r is 0 or 1;  
s is 0 or 1;

R<sup>1</sup> is selected from:

- 1) —(C=O)C<sub>1</sub>-C<sub>10</sub>-alkyl,
- 2) —(C=O)aryl,
- 3) —(C=O)C<sub>3</sub>-C<sub>8</sub>-cycloalkyl,
- 4) —(C=O)heterocyclyl,
- 5) (C=O)NR<sup>c</sup>R<sup>c'</sup>,
- 6) —(C=S)NR<sup>e</sup>R<sup>e'</sup>,
- 7) —SO<sub>2</sub>NR<sup>e</sup>R<sup>e'</sup>,
- 8) —SO<sub>2</sub>C<sub>1</sub>-C<sub>10</sub>-alkyl,
- 9) —SO<sub>2</sub>-aryl, and
- 10) —SO<sub>2</sub>-heterocyclyl,

said alkyl, aryl, cycloalkyl, and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;

R<sup>2</sup> is selected from:

- 1) aryl,
- 2) —C<sub>1</sub>-C<sub>6</sub>-aralkyl,
- 3) —C<sub>3</sub>-C<sub>8</sub>-cycloalkyl, and
- 4) —heterocyclyl,

said aryl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;

R<sup>3</sup>, R<sup>4</sup> and R<sup>8</sup> are independently selected from:

- 1) H, and
- 2) C<sub>1</sub>-C<sub>10</sub> alkyl, and
- 3) —C<sub>1</sub>-C<sub>6</sub>-perfluoroalkyl,

~~said alkyl, aryl, alkenyl, alkynyl, cycloalkyl, aralkyl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;~~

R<sup>6</sup> is selected from:

- 1) aryl,
- 2) ~~C<sub>1</sub>-C<sub>6</sub> aralkyl,~~
- 3) ~~C<sub>3</sub>-C<sub>8</sub> cycloalkyl, and~~
- 4) ~~heterocyclyl,~~

said aryl, ~~cycloalkyl, aralkyl and heterocyclyl~~ is optionally substituted with one or more substituents selected from R<sup>10</sup>;

R<sup>10</sup> is independently selected from:

- 1) (C=O)<sub>a</sub>O<sub>b</sub>C<sub>1</sub>-C<sub>10</sub> alkyl,
- 2) (C=O)<sub>a</sub>O<sub>b</sub>aryl,
- 3) C<sub>2</sub>-C<sub>10</sub> alkenyl,
- 4) C<sub>2</sub>-C<sub>10</sub> alkynyl,
- 5) (C=O)<sub>a</sub>O<sub>b</sub> heterocyclyl,
- 6) CO<sub>2</sub>H,
- 7) halo,
- 8) CN,
- 9) OH,
- 10) O<sub>b</sub>C<sub>1</sub>-C<sub>6</sub> perfluoroalkyl,
- 11) O<sub>a</sub>(C=O)<sub>b</sub>NR<sup>12</sup>R<sup>13</sup>,
- 12) S(O)<sub>m</sub>R<sup>a</sup>,
- 13) S(O)<sub>2</sub>NR<sup>12</sup>R<sup>13</sup>,
- 14) ~~oxo,~~
- 15) CHO, and
- 16) ~~(N=O)R<sup>12</sup>R<sup>13</sup>, or~~
- 17) (C=O)<sub>a</sub>O<sub>b</sub>C<sub>3</sub>-C<sub>8</sub> cycloalkyl,

said alkyl, aryl, alkenyl, alkynyl, heterocyclyl, and cycloalkyl optionally substituted with one, two or three substituents selected from R<sup>11</sup>;

R<sup>11</sup> is selected from:

- 1) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>1</sub>-C<sub>10</sub>)alkyl,
- 2) O<sub>r</sub>(C<sub>1</sub>-C<sub>3</sub>)perfluoroalkyl,
- 3) oxo,
- 4) OH,
- 5) halo,
- 6) CN,
- 7) (C<sub>2</sub>-C<sub>10</sub>)alkenyl,
- 8) (C<sub>2</sub>-C<sub>10</sub>)alkynyl,
- 9) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>3</sub>-C<sub>6</sub>)cycloalkyl,
- 10) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>0</sub>-C<sub>6</sub>)alkylene-aryl,
- 11) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>0</sub>-C<sub>6</sub>)alkylene-heterocyclyl,
- 12) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>0</sub>-C<sub>6</sub>)alkylene-N(R<sup>b</sup>)<sub>2</sub>,
- 13) C(O)R<sup>a</sup>,
- 14) (C<sub>0</sub>-C<sub>6</sub>)alkylene-CO<sub>2</sub>R<sup>a</sup>,
- 15) C(O)H,
- 16) (C<sub>0</sub>-C<sub>6</sub>)alkylene-CO<sub>2</sub>H,
- 17) C(O)N(R<sup>b</sup>)<sub>2</sub>,
- 18) S(O)<sub>m</sub>R<sup>a</sup>, and
- 19) S(O)<sub>2</sub>N(R<sup>b</sup>)<sub>2</sub>;

said alkyl, alkenyl, alkynyl, cycloalkyl, aryl, and heterocyclyl is optionally substituted with up to three substituents selected from R<sup>b</sup>, OH, (C<sub>1</sub>-C<sub>6</sub>)alkoxy, halogen, CO<sub>2</sub>H, CN, O(C=O)C<sub>1</sub>-C<sub>6</sub> alkyl, oxo, and N(R<sup>b</sup>)<sub>2</sub>;

R<sup>12</sup> and R<sup>13</sup> are independently selected from:

- 1) H,
- 2) (C=O)O<sub>b</sub>C<sub>1</sub>-C<sub>10</sub> alkyl,
- 3) (C=O)O<sub>b</sub>C<sub>3</sub>-C<sub>8</sub> cycloalkyl,
- 4) (C=O)O<sub>b</sub>aryl,
- 5) (C=O)O<sub>b</sub>heterocyclyl,
- 6) C<sub>1</sub>-C<sub>10</sub> alkyl,

- 7) aryl,
- 8) C<sub>2</sub>-C<sub>10</sub> alkenyl,
- 9) C<sub>2</sub>-C<sub>10</sub> alkynyl,
- 10) heterocyclyl,
- 11) C<sub>3</sub>-C<sub>8</sub> cycloalkyl,
- 12) SO<sub>2</sub>R<sup>a</sup>, and
- 13) (C=O)NR<sup>b</sup><sub>2</sub>,

said alkyl, cycloalkyl, aryl, heterocyclyl, alkenyl, and alkynyl is optionally substituted with one, two or three substituents selected from R<sup>11</sup>, or

R<sup>12</sup> and R<sup>13</sup> can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 5-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R<sup>11</sup>;

R<sup>a</sup> is independently selected from: (C<sub>1</sub>-C<sub>6</sub>)alkyl, (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, aryl, and heterocyclyl;

R<sup>b</sup> is independently selected from: H, (C<sub>1</sub>-C<sub>6</sub>)alkyl, aryl, heterocyclyl, (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, (C=O)OC<sub>1</sub>-C<sub>6</sub> alkyl, (C=O)C<sub>1</sub>-C<sub>6</sub> alkyl or S(O)<sub>2</sub>R<sup>a</sup>; and

R<sup>c</sup> and R<sup>c'</sup> are independently selected from: H, (C<sub>1</sub>-C<sub>6</sub>)alkyl, aryl, heterocyclyl and (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl or

R<sup>c</sup> and R<sup>c'</sup> can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 5-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R<sup>11</sup>.

7. (Currently amended) The compound according to Claim 6 or a pharmaceutically acceptable salt or stereoisomer thereof, wherein:

R<sup>1</sup> is selected from:

- 1) (C=O)NR<sup>c</sup>R<sup>c'</sup>,
- 2) —SO<sub>2</sub>NR<sup>e</sup>R<sup>e'</sup>, and
- 3) —SO<sub>2</sub>C<sub>1</sub>-C<sub>10</sub> alkyl,

~~said alkyl is optionally substituted with one, two or three substituents selected from R<sup>10</sup>;~~

R<sup>2</sup> is selected from:

- 1) aryl, and
- 2) ~~heteroaryl,~~

~~said aryl and heteroaryl is optionally substituted with one or more substituents selected from R<sup>10</sup>;~~

R<sup>3</sup>, R<sup>4</sup> and R<sup>8</sup> are independently selected from:

- 1) H, and
- 2) C<sub>1</sub>-C<sub>10</sub> alkyl,

~~said alkyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;~~

R<sup>6</sup> is selected from:

- 1) aryl, and
- ~~2) heterocyclyl,~~

~~said alkyl, aryl and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>; and~~

~~R<sup>10</sup>, R<sup>11</sup>, R<sup>12</sup>, R<sup>13</sup>, R<sup>a</sup>, R<sup>b</sup>, R<sup>e</sup> and R<sup>e'</sup> are as described immediately above.~~

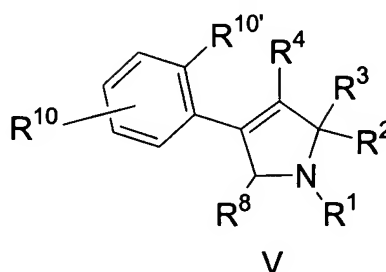
8. (Currently amended) The compound according to Claim 7 or a pharmaceutically acceptable salt or stereoisomer thereof, wherein:

~~R<sup>2</sup> and R<sup>6</sup> are independently selected from~~ are phenyl ~~or pyridyl~~, optionally substituted with one or two substituents selected from R<sup>10</sup>; ~~and~~

~~R<sup>1</sup>, R<sup>3</sup>, R<sup>4</sup> and R<sup>8</sup> are as described in Claim 7.~~

9. (Cancelled)

10. (Currently amended) The compound according to Claim 5 of the Formula V,



wherein:

a is 0 or 1;  
b is 0 or 1;  
m is 0, 1, or 2;  
r is 0 or 1;  
s is 0 or 1;

R<sup>1</sup> is selected from:

- 1) —(C=O)C<sub>1</sub>-C<sub>10</sub>-alkyl,
- 2) —(C=O)aryl,
- 3) —(C=O)C<sub>3</sub>-C<sub>8</sub>-cycloalkyl,
- 4) —(C=O)heterocyclyl,
- 5) (C=O)NR<sup>c</sup>R<sup>c'</sup>,
- 6) —(C=S)NR<sup>e</sup>R<sup>e'</sup>,
- 7) —SO<sub>2</sub>NR<sup>e</sup>R<sup>e'</sup>,
- 8) —SO<sub>2</sub>C<sub>1</sub>-C<sub>10</sub>-alkyl,
- 9) —SO<sub>2</sub>-aryl, and
- 10) —SO<sub>2</sub>-heterocyclyl,

said-alkyl, aryl, cycloalkyl, and heterocyclyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;



R<sup>2</sup> is selected from:

- 1) aryl,
- 2) ~~C<sub>1</sub>-C<sub>6</sub> alkyl,~~
- 3) ~~C<sub>3</sub>-C<sub>8</sub> cycloalkyl, and~~
- 4) ~~heterocyclyl,~~

said aryl, ~~cycloalkyl, alkyl and heterocyclyl~~ is optionally substituted with one or more substituents selected from R<sup>10</sup>;

R<sup>3</sup>, R<sup>4</sup> and R<sup>8</sup> are independently selected from:

- 1) H, and
- 2) C<sub>1</sub>-C<sub>10</sub> alkyl, and
- 3) ~~C<sub>1</sub>-C<sub>6</sub> perfluoroalkyl,~~

~~said alkyl is optionally substituted with one or more substituents selected from R<sup>10</sup>;~~

R<sup>10</sup> is independently selected from:

- 1) (C=O)<sub>a</sub>O<sub>b</sub>C<sub>1</sub>-C<sub>10</sub> alkyl,
- 2) (C=O)<sub>a</sub>O<sub>b</sub>aryl,
- 3) C<sub>2</sub>-C<sub>10</sub> alkenyl,
- 4) C<sub>2</sub>-C<sub>10</sub> alkynyl,
- 5) (C=O)<sub>a</sub>O<sub>b</sub> heterocyclyl,
- 6) CO<sub>2</sub>H,
- 7) halo,
- 8) CN,
- 9) OH,
- 10) O<sub>b</sub>C<sub>1</sub>-C<sub>6</sub> perfluoroalkyl,
- 11) O<sub>a</sub>(C=O)<sub>b</sub>NR<sup>12</sup>R<sup>13</sup>,
- 12) S(O)<sub>m</sub>R<sup>a</sup>,
- 13) S(O)<sub>2</sub>NR<sup>12</sup>R<sup>13</sup>,
- 14) ~~oxo,~~
- 15) CHO, and
- 16) ~~(N=O)R<sup>12</sup>R<sup>13</sup>, or~~

17)  $(\text{C}=\text{O})_a\text{O}_b\text{C}_3\text{-C}_8$  cycloalkyl,  
said alkyl, aryl, alkenyl, alkynyl, heterocyclyl, and cycloalkyl optionally substituted with one,  
two or three substituents selected from  $\text{R}^{11}$ ;  
 $\text{R}^{10'}$  is halogen;

$\text{R}^{11}$  is selected from:

- 1)  $(\text{C}=\text{O})_r\text{O}_s(\text{C}_1\text{-C}_{10})$ alkyl,
- 2)  $\text{O}_r(\text{C}_1\text{-C}_3)$ perfluoroalkyl,
- 3) oxo,
- 4) OH,
- 5) halo,
- 6) CN,
- 7)  $(\text{C}_2\text{-C}_{10})$ alkenyl,
- 8)  $(\text{C}_2\text{-C}_{10})$ alkynyl,
- 9)  $(\text{C}=\text{O})_r\text{O}_s(\text{C}_3\text{-C}_6)$ cycloalkyl,
- 10)  $(\text{C}=\text{O})_r\text{O}_s(\text{C}_0\text{-C}_6)$ alkylene-aryl,
- 11)  $(\text{C}=\text{O})_r\text{O}_s(\text{C}_0\text{-C}_6)$ alkylene-heterocyclyl,
- 12)  $(\text{C}=\text{O})_r\text{O}_s(\text{C}_0\text{-C}_6)$ alkylene- $\text{N}(\text{R}^b)_2$ ,
- 13)  $\text{C}(\text{O})\text{R}^a$ ,
- 14)  $(\text{C}_0\text{-C}_6)$ alkylene- $\text{CO}_2\text{R}^a$ ,
- 15)  $\text{C}(\text{O})\text{H}$ ,
- 16)  $(\text{C}_0\text{-C}_6)$ alkylene- $\text{CO}_2\text{H}$ ,
- 17)  $\text{C}(\text{O})\text{N}(\text{R}^b)_2$ ,
- 18)  $\text{S}(\text{O})_m\text{R}^a$ , and
- 19)  $\text{S}(\text{O})_2\text{N}(\text{R}^b)_2$ ;

said alkyl, alkenyl, alkynyl, cycloalkyl, aryl, alkylene and heterocyclyl is optionally substituted with up to three substituents selected from  $\text{R}^b$ , OH,  $(\text{C}_1\text{-C}_6)$ alkoxy, halogen,  $\text{CO}_2\text{H}$ , CN,  $\text{O}(\text{C}=\text{O})\text{C}_1\text{-C}_6$  alkyl, oxo, and  $\text{N}(\text{R}^b)_2$ ;

$\text{R}^{12}$  and  $\text{R}^{13}$  are independently selected from:

- 1) H,
- 2)  $(\text{C}=\text{O})\text{O}_b\text{C}_1\text{-C}_{10}$  alkyl,

- 3)  $(\text{C}=\text{O})\text{O}_b\text{C}_3\text{-C}_8$  cycloalkyl,
- 4)  $(\text{C}=\text{O})\text{O}_b$ aryl,
- 5)  $(\text{C}=\text{O})\text{O}_b$ heterocyclyl,
- 6)  $\text{C}_1\text{-C}_{10}$  alkyl,
- 7) aryl,
- 8)  $\text{C}_2\text{-C}_{10}$  alkenyl,
- 9)  $\text{C}_2\text{-C}_{10}$  alkynyl,
- 10) heterocyclyl,
- 11)  $\text{C}_3\text{-C}_8$  cycloalkyl,
- 12)  $\text{SO}_2\text{R}^a$ , and
- 13)  $(\text{C}=\text{O})\text{NR}^b_2$ ,

said alkyl, cycloalkyl, aryl, heterocyclyl, alkenyl, and alkynyl is optionally substituted with one, two or three substituents selected from  $\text{R}^{11}$ , or

$\text{R}^{12}$  and  $\text{R}^{13}$  can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 5-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from  $\text{R}^{11}$ ;

$\text{R}^a$  is independently selected from:  $(\text{C}_1\text{-C}_6)$ alkyl,  $(\text{C}_3\text{-C}_6)$ cycloalkyl, aryl, and heterocyclyl;

$\text{R}^b$  is independently selected from: H,  $(\text{C}_1\text{-C}_6)$ alkyl, aryl, heterocyclyl,  $(\text{C}_3\text{-C}_6)$ cycloalkyl,  $(\text{C}=\text{O})\text{OC}_1\text{-C}_6$  alkyl,  $(\text{C}=\text{O})\text{C}_1\text{-C}_6$  alkyl or  $\text{S}(\text{O})_2\text{R}^a$ ; and

$\text{R}^c$  and  $\text{R}^{c'}$  are independently selected from: H,  $(\text{C}_1\text{-C}_6)$ alkyl, aryl, heterocyclyl and  $(\text{C}_3\text{-C}_6)$ cycloalkyl or

$\text{R}^c$  and  $\text{R}^{c'}$  can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 5-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said

monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R<sup>11</sup>.

11. (Currently amended) A compound selected from:

4-(2-chloro-5-fluorophenyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(+)-4-(2,5-difluorophenyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(-)-4-(2,5-difluorophenyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(5-chloro-2-fluorophenyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2-fluorophenyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2-fluoro-5-methylphenyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(5-bromo-2-fluorophenyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4- {[4-(5-chloro-2-fluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}morpholine;

4- {[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}morpholine;

N,N-dimethyl-2,4-diphenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

~~3-[2-fluoro-5-(trifluoromethyl)phenyl]-N,N-dimethyl-5-phenyl-2,3-dihydro-1H-pyrrole-1-carboxamide;~~

2-(3-fluorophenyl)-4-(2,5-difluorophenyl)-N,N-dimethyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(4-fluorophenyl)-N,N-dimethyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(2-fluorophenyl)-N,N-dimethyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

2-(3-bromophenyl)-4-(2,5-difluorophenyl)- N,N-dimethyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

2-(3-aminophenyl)-4-(2,5-difluorophenyl)- N,N-dimethyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-methylphenyl)-N,N-dimethyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(5-chloro-2-fluorophenyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

~~4-(2,5-difluorophenyl)-1-(methylsulfonyl)-2-phenyl-2,5-dihydro-1H-pyrrole;~~

~~4-(2,5-difluorophenyl)-1-(ethylsulfonyl)-2-phenyl-2,5-dihydro-1H-pyrrole;~~

~~4-(2,5-difluorophenyl)-2-phenyl-1-(propylsulfonyl)-2,5-dihydro-1H-pyrrole;~~

~~4-(2,5-difluorophenyl)-1-(isopropylsulfonyl)-2-phenyl-2,5-dihydro-1H-pyrrole;~~

~~4-(5-chloro-2-fluorophenyl)-1-(methylsulfonyl)-2-phenyl-2,5-dihydro-1H-pyrrole;~~

~~4-(5-chloro-2-fluorophenyl)-1-(isopropylsulfonyl)-2-phenyl-2,5-dihydro-1H-pyrrole;~~

~~4-(2-fluoro-5-methylphenyl)-1-(isopropylsulfonyl)-2-phenyl-2,5-dihydro-1H-pyrrole;~~

~~2-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]sulfonyl]ethanamine;~~

~~2-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]sulfonyl]-N,N-dimethylethanamine;~~

~~1-acetyl-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrole;~~

~~4-(2-chloro-5-fluorophenyl)-1-pivaloyl-2-phenyl-2,5-dihydro-1H-pyrrole;~~

~~4-(2-chloro-5-fluorophenyl)-1-isobutyryl-2-phenyl-2,5-dihydro-1H-pyrrole;~~

~~4-(2,5-difluorophenyl)-1-(2,2-dimethylpropanoyl)-2-phenyl-2,5-dihydro-1H-pyrrole;~~

~~1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-methyl-1-oxopropan-2-ol;~~

~~1-[4-(5-chloro-2-fluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-methyl-1-oxopropan-2-ol;~~

~~1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-methyl-1-oxopropan-2-amine;~~

~~4-(2-fluoro-5-isocyanophenyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;~~

~~4-(2,5-difluorophenyl)-2-phenyl-1-(trifluoroacetyl)-2,5-dihydro-1H-pyrrole;~~

~~4-(5-chloro-2-fluorophenyl)-2-phenyl-1-(trifluoroacetyl)-2,5-dihydro-1H-pyrrole;~~

~~(1S)-1-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-2-methylpropylamine;~~

~~(1R)-1-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-2-methylpropylamine;~~

~~4-(2,5-difluorophenyl)-2-phenyl-1-L-prolyl-2,5-dihydro-1H-pyrrole;~~

~~4-(2,5-difluorophenyl)-2-phenyl-1-D-prolyl-2,5-dihydro-1H-pyrrole;~~

~~(4R)-4-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-1,3-thiazolidine;~~

~~methyl-(3S)-3-amino-4-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-4-oxobutanoate;~~

~~(4S)-4-amino-5-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-5-oxopentanamide;~~

~~(1S)-1-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-3-(methylthio)propylamine;~~

~~(1S)-1-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-3-(methylsulfonyl)propylamine;~~

~~(2S)-2-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]piperidine;~~

~~(1S)-1-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]pentylamine;~~

~~(1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxo-1-(thien-2-ylmethyl)ethylamine;~~

~~4-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-1,1-dioxidotetrahydro-2H-thiopyran-4-ylamine;~~

~~(2S)-1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-methyl-1-oxopropan-2-amine;~~

~~(1S)-1-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]propylamine;~~

~~(1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxo-1-phenylethanamine;~~

~~(1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxo-1-phenylethanamine;~~

~~(4S)-4-amino-5-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-5-oxopentanamide~~

~~3-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3-oxopropan-1-amine;~~

~~(1S,2S)-1-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-2-methylbutylamine;~~

~~(1S)-1-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]butylamine;~~

~~(1S)-1-cyclopropyl-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethanamine;~~

~~1-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]cyclopropanamine;~~

~~1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-oxopropan-2-amine;~~

~~(1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-methyl-2-oxoethylamine;~~

~~(1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxo-1-(pyridin-2-ylmethyl)ethylamine;~~

~~(1S)-1-cyclohexyl-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethanamine;~~

~~(1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-(4-iodobenzyl)-2-oxoethylamine;~~

~~(1S)-1-benzyl-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethylamine;~~

~~4-((2S)-2-amino-3-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3-oxopropyl)phenol;~~



~~(3S)-3-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-1,2,3,4-tetrahydroisoquinoline;~~

~~(1S)-1-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-3-phenylpropylamine;~~

~~(1S)-1-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-3-methylbutylamine;~~

~~(1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxo-1-(pyridin-3-ylmethyl)ethylamine;~~

~~1-[(2S)-azetidin-2-ylcarbonyl]-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrole;~~

~~(3S)-3-amino-4-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-4-oxobutanamide;~~

~~4-(2,5-difluorophenyl)-1-[(2S)-2,5-dihydro-1H-pyrrol-2-ylcarbonyl]-2-phenyl-2,5-dihydro-1H-pyrrole;~~

~~4-(2,5-difluorophenyl)-1-[(2-methylazetidin-2-yl)carbonyl]-2-phenyl-2,5-dihydro-1H-pyrrole;~~

~~(1S)-1-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-2,2-dimethylpropylamine;~~

~~methyl (4S)-4-amino-5-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-5-oxopentanoate;~~

~~4-(2,5-difluorophenyl)-2-phenyl-1-[(2S,3S)-2-phenylpyrrolidin-3-yl]carbonyl]-2,5-dihydro-1H-pyrrole;~~

~~4-(2,5-difluorophenyl)-2-phenyl-1-[(5-phenylpyrrolidin-3-yl)carbonyl]-2,5-dihydro-1H-pyrrole;~~

~~(2S)-2-amino-3-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3-oxopropan-1-ol;~~

~~(2R,3S)-3-amino-4-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-4-oxobutan-2-ol;~~

~~(1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-(methoxymethyl)-2-oxoethylamine;~~

~~4-(2,5-difluorophenyl)-2-phenyl-1-(pyrrolidin-3-ylcarbonyl)-2,5-dihydro-1H-pyrrole;~~

~~4-(2,5-difluorophenyl)-2-phenyl-1-[(3-phenylpyrrolidin-3-yl)acetyl]-2,5-dihydro-1H-pyrrole;~~

~~(1S)-1-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-3,3-difluoropropylamine;~~

~~(1S)-3-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3-oxo-1-phenylpropan-1-amine;~~

~~4-(2,5-difluorophenyl)-2-phenyl-1-[(4S)-4-phenyl-L-prolyl]-2,5-dihydro-1H-pyrrole;~~

~~1-{2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}cyclohexanamine;~~

~~2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethanamine;~~

~~4-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]piperidin-4-amine;~~

~~(1S,3R)-3-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]cyclopentanamine;~~

~~(1R,4S)-4-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]cyclopent-2-en-1-amine;~~

~~(1S,4R)-4-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]cyclopent-2-en-1-amine;~~

~~(1S)-1-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]but-3-ynylamine;~~

~~(1R)-3-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3-oxo-1-phenylpropan-1-amine;~~

~~3-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-2-phenylpiperidine;~~

~~(1S)-1-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]but-3-enylamine;~~

~~(2S)-3-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-(methylamino)-3-oxopropan-1-ol;~~

~~(3R,5S)-5-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]pyrrolidin-3-ol;~~

~~(1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxo-1-(1,3-thiazol-4-ylmethyl)ethylamine;~~

~~(1R)-1-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]but-3-enylamine;~~

~~(2S)-1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N,3-dimethyl-1-oxobutan-2-amine;~~

~~(2S)-1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N,4-dimethyl-1-oxopentan-2-amine;~~

~~(1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-[(1-methyl-1H-imidazol-4-yl)methyl]-2-oxoethylamine;~~

~~4-(2,5-difluorophenyl)-1-(N-6-formyl-L-lysyl)-2-phenyl-2,5-dihydro-1H-pyrrole;~~

~~(2S,3S)-1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N,3-dimethyl-1-oxopentan-2-amine;~~

~~(1S)-1-(cyclohexylmethyl)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethylamine;~~

~~(1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-(1H-indol-3-ylmethyl)-2-oxoethylamine;~~

~~(1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-(isocyanomethyl)-2-oxoethylamine;~~

~~(1S)-1-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-3,3-dimethylbutylamine;~~

~~1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2,3-dimethyl-1-oxobutan-2-amine;~~

~~1-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]cyclohexanamine;~~

~~1-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]cyclopentanamine;~~

~~(1S)-3-(benzyloxy)-1-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]propylamine;~~

~~1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2,3-dimethyl-1-oxobutan-2-amine;~~

~~1-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]cyclopent-3-en-1-amine;~~

~~(1S)-1-cyclopentyl-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethanamine;~~

~~4-(2,5-difluorophenyl)-1-(2-methylpropyl)-2-phenyl-2,5-dihydro-1H-pyrrole;~~

~~1-[4-(5-chloro-2-fluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-methyl-1-oxopropan-2-amine;~~

~~(1S)-1-[[4-(5-chloro-2-fluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-2-methylpropylamine;~~

~~(1S)-2-[4-(5-chloro-2-fluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-cyclopropyl-2-oxoethanamine;~~

~~(1S,2S)-1-[[4-(5-chloro-2-fluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-2-methylbutylamine;~~

~~(1S)-1-[[4-(5-chloro-2-fluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]pentylamine;~~

~~(1S)-1-[[4-(5-chloro-2-fluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-3,3-dimethylbutylamine;~~

~~(1S)-1-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-2,2-dimethylpropylamine;~~

~~(1S)-1-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-2-methylpropylamine;~~

~~(1S)-1-cyclohexyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethanamine;~~

~~(1S)-1-[[[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]but-3-enylamine;~~

~~(1S)-1-[[[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]but-3-ynylamine;~~

~~(1S)-1-benzyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethylamine;~~

~~(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethanamine;~~

~~1-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-methyl-1-oxopropan-2-amine;~~

~~(1S)-1-[[[(2S)-4-(5-chloro-2-fluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-2,2-dimethylpropylamine;~~

~~(1S)-1-[[[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]pentylamine;~~

~~(1S)-1-[[[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-3-methylbutylamine;~~

~~(1S)-1-[[[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-3,3-dimethylbutylamine;~~

~~1-cyclopropyl-3-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3-oxopropan-1-amine;~~

~~(1S)-2-[(2S)-4-(5-chloro-2-fluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-cyclopropyl-2-oxoethanamine;~~

~~(1S)-1-[[[(2S)-4-(5-chloro-2-fluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-2-methylpropylamine;~~

~~(1S,2S)-1-[[[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-2-methylbutylamine;~~

~~4-(5-chloro-2-fluorophenyl)-2-(3-hydroxyphenyl)-1-(2-methylalanyl)-2,5-dihydro-1H-pyrrole;~~

~~(2S)-4-(5-chloro-2-fluorophenyl)-2-(3-hydroxyphenyl)-1-(2-methylalanyl)-2,5-dihydro-1H-pyrrole;~~

~~4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-1-L-valyl-2,5-dihydro-1H-pyrrole;~~

~~4-(5-chloro-2-fluorophenyl)-2-(3-hydroxyphenyl)-1-L-valyl-2,5-dihydro-1H-pyrrole;~~

~~(2S)-4-(5-chloro-2-fluorophenyl)-2-(3-hydroxyphenyl)-1-L-valyl-2,5-dihydro-1H-pyrrole;~~

~~4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-1-(2-methylalanyl)-2,5-dihydro-1H-pyrrole;~~

~~3-[1-[(2S)-2-amino-2-cyclopropylethanoyl]-4-(5-chloro-2-fluorophenyl)-2,5-dihydro-1H-pyrrol-2-yl]phenol;~~

~~4-(5-chloro-2-fluorophenyl)-2-(3-hydroxyphenyl)-1-L-isoleucyl-2,5-dihydro-1H-pyrrole;~~

~~4-(5-chloro-2-fluorophenyl)-2-(3-hydroxyphenyl)-1-L-norleucyl-2,5-dihydro-1H-pyrrole;~~

~~(2S)-4-(5-chloro-2-fluorophenyl)-2-(3-hydroxyphenyl)-1-(3-methyl-L-valyl)-2,5-dihydro-1H-pyrrole;~~

~~(2S)-4-(2,5-Difluorophenyl)-N-methyl-2-phenyl-N-piperidin-4-yl-2,5-dihydro-1H-pyrrole-1-carboxamide;~~

(2S)-4-(2,5-Difluorophenyl)-N-methyl-2-phenyl-N-(piperidin-4-ylmethyl)-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(5-Chloro-2-fluorophenyl)-N-methyl-2-phenyl-N-[(3R)-pyrrolidin-3-yl]-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(5-chloro-2-fluorophenyl)-N-methyl-2-phenyl-N-piperidin-4-yl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(5-chloro-2-fluorophenyl)-N-methyl-2-phenyl-N-[(3S)-pyrrolidin-3-yl]-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(2,5-difluorophenyl)-N-methyl-2-phenyl-N-pyrrolidin-3-yl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(2,5-difluorophenyl)-N-methyl-N-[(3S)-1-methylpyrrolidin-3-yl]-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(2,5-difluorophenyl)-N-methyl-N-[(3R)-1-methylpyrrolidin-3-yl]-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(2,5-difluorophenyl)-N-methyl-N-[(1-methyl-5-oxopyrrolidin-2-yl)methyl]-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(2,5-difluorophenyl)-N-(1,3-dioxolan-2-ylmethyl)-N-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(2,5-difluorophenyl)-N-methyl-2-phenyl-N-tetrahydrofuran-3-yl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-N-(1-allylpiperidin-4-yl)-4-(2,5-difluorophenyl)-N-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;



allyl 4-[[[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl](methyl)amino]piperidine-1-carboxylate;

allyl 4-[[[[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl](methyl)amino]methyl]piperidine-1-carboxylate;

(2S)-4-(2,5-difluorophenyl)-N-methyl-N-(1-methylpiperidin-4-yl)-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-Difluorophenyl)-N-methyl-N-[(1-methylpiperidin-3-yl)methyl]-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-N-methyl-2-phenyl-N-(pyridin-3-ylmethyl)-2,5-dihydro-1H-pyrrole-1-carboxamide;

N-benzyl-4-(2,5-difluorophenyl)-N-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-N-methyl-N-[(1-methyl-1H-pyrazol-4-yl)methyl]-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-N-[2-(dimethylamino)ethyl]-N-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-N-(2-hydroxyethyl)-N-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-N-isobutyl-N-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-N-methyl-2-phenyl-N-(2-pyridin-2-ylethyl)-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-N-(2-methoxyethyl)-N-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-N-(2,3-dihydroxypropyl)-N-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-N-methyl-2-phenyl-N-(2-phenylethyl)-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-phenyl-N-propyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(5-chloro-2-fluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-piperidin-4-yl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-piperidin-4-yl-2,5-dihydro-1H-pyrrole-1-carboxamide;

~~1-Acetyl 4-(2,5-difluorophenyl)-2-methyl-2-phenyl-2,5-dihydro-1H-pyrrole;~~

~~(2S)-1-[4-(2,5-difluorophenyl)-2-methyl-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3-methyl-1-oxobutan-2-amine;~~

(2S)-4-(2,5-difluorophenyl)-N,N,2-trimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(5-chloro-2-fluorophenyl)-N,N,2-trimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(5-chloro-2-fluorophenyl)-2-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(5-chloro-2-fluorophenyl)-N-ethyl-2-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

~~(2S)-1-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3,3-dimethyl-1-oxobutan-2-ol;~~

~~(2S)-1-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3-methyl-1-oxobutan-2-ol;~~

~~(2S,3S)-1-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3-methyl-1-oxopentan-2-ol;~~

~~1-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-oxohexan-2-ol;~~

~~(2S)-1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-oxo-3-phenylpropan-2-ol;~~

~~(2S)-1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-4-methyl-1-oxopentan-2-ol;~~

~~(1S)-1-cyclohexyl-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethanol;~~

~~(2S)-1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3,3-dimethyl-1-oxobutan-2-ol;~~

~~N-1-[(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl]-N-2,N-2-dimethylglycinamide;~~

~~N-1-[(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl]-N-2-methylglycinamide;~~

~~N-1-[(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl]glycinamide;~~

~~N-1-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-2-methylalaninamide;~~

~~N-1-((1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)glycinamide;~~

~~N-1-((1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-N-2,N-2-dimethylglycinamide;~~

~~N-1-((1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-N-2,N-2-dimethylglycinamide, N-oxide;~~

~~N-1-((1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-2-methylalaninamide;~~

~~N-1-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-N-2,N-2-dimethylglycinamide n-oxide;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-2-pyrrolidin-1-ylacetamide;~~

~~2-azetidin-1-yl-N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)acetamide;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-2-morpholin-4-ylacetamide;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-2-piperazin-1-ylacetamide;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-2-(4-methylpiperazin-1-yl)acetamide;~~

~~2-azetidin-1-yl-N-((1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)acetamide;~~

~~N-((1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-2-pyrrolidin-1-ylacetamide;~~

~~N-((1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-2-piperidin-1-ylacetamide;~~

~~N-((1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-2-morpholin-4-ylacetamide;~~

~~N-1-((1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-N-2-(2-hydroxyethyl)glycinamide;~~

~~N-((1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-2-(4-methylpiperazin-1-yl)acetamide;~~

~~N-1-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-N-2-isopropylglycinamide;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)acetamide;~~

~~N-1-((1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-N-2-ethylglycinamide;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-2-hydroxyacetamide;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)piperazine-1-carboxamide;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-N'-piperidin-4-ylurea;~~

~~4-amino-N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)piperidine-1-carboxamide;~~

~~N-(2-aminoethyl)-N'-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)urea;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-N'-(3-morpholin-4-ylpropyl)urea;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-N'-[2-(dimethylamino)ethyl]urea;~~

~~2-azetidin-1-yl-N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)ethanesulfonamide~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-2-(isopropylamino)ethanesulfonamide;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-2-pyrrolidin-1-ylethanesulfonamide;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-2-morpholin-4-ylethanesulfonamide;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-2-piperazin-1-ylethanesulfonamide;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-2-(4-methylpiperazin-1-yl)ethanesulfonamide;~~

N-(tert-butyl)-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]acetamide;  
2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-isopropylacetamide;  
(2S)-1-(2-azetidin-1-yl-2-oxoethyl)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrole;  
(2S)-4-(2,5-difluorophenyl)-1-(2-oxo-2-pyrrolidin-1-ylethyl)-2-phenyl-2,5-dihydro-1H-pyrrole;  
4-[[[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]acetyl}morpholine;  
1-[[[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]acetyl}piperazine;  
1-[[[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]acetyl}-4-methylpiperazine;  
2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-isopropylbutanamide;  
4-{2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]butanoyl}morpholine;  
2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-ethylacetamide;  
N-cyclobutyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]acetamide;  
2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-ethylpropanamide;  
N-cyclobutyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]propanamide;  
2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-methylpropanamide;  
2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-isopropylpropanamide;  
N-(tert-butyl)-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]propanamide;

4-{2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]propanoyl}morpholine;

~~(3S)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-ethyl-2,2-dimethyl-4-oxobutanamide;~~

~~(3S)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2,2-dimethyl-4-oxo-N-piperidin-4-ylbutanamide;~~

~~(3S)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2,2-dimethyl-4-oxobutanoic acid;~~

~~(3S)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N,N,2,2-tetramethyl-4-oxobutanamide;~~

~~(1S)-1-[[[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-2,2-dimethyl-3-oxo-3-piperazin-1-yl]propylamine;~~

~~(3S)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-isopropyl-2,2-dimethyl-4-oxobutanamide;~~

~~(3S)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N,2,2-trimethyl-4-oxobutanamide;~~

~~(3R)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N,N,2,2-tetramethyl-4-oxobutanamide;~~

~~(3R)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2,2-dimethyl-4-oxobutanoic acid;~~

~~(1R)-1-[[[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-2,2-dimethyl-3-oxo-3-piperazin-1-yl]propylamine;~~



~~2-(((1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}amino)-N-ethylacetamide;~~

~~2-(((1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}amino)-N-methylacetamide;~~

~~2-(((1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}amino)-N,N-dimethylacetamide;~~

~~2-(((1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}amino)-N-methyl-N-ethylacetamide;~~

~~2-(((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}amino)-N-methylacetamide;~~

~~2-(((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}amino)-N-ethylacetamide;~~

~~2-(((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}amino)-N,N-dimethylacetamide;~~

~~2-(((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}amino)-N-isopropylacetamide;~~

~~2-(((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}amino)-N-ethyl-N-methylacetamide;~~

~~2-(((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}amino)-N,N-diethylacetamide;~~

~~(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxo-N-(2-oxo-2-pyrrolidin-1-ylethyl)ethanamine;~~

~~(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-(2-morpholin-4-yl-2-oxoethyl)-2-oxoethanamine;~~

~~1-[(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl]amino)acetyl]piperidin-4-ol;~~

~~(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-[2-(4-methylpiperazin-1-yl)-2-oxoethyl]-2-oxoethanamine;~~

~~(1S)-N-(2-azetidin-1-yl-2-oxoethyl)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethanamine;~~

~~(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-[2-(1,1-dioxidothiomorpholin-4-yl)-2-oxoethyl]-2-oxoethanamine;~~

~~(1S)-N-[2-(4-acetylpiperazin-1-yl)-2-oxoethyl]-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethanamine;~~

~~(1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-(2-morpholin-4-yl-2-oxoethyl)-2-oxoethanamine;~~

~~(1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxo-N-(2-oxo-2-pyrrolidin-1-ylethyl)ethanamine;~~

~~2-((1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)amino)-N-isopropylacetamide;~~

~~2-(dimethylamino)ethyl (1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethylcarbamate;~~

~~1-methylpiperidin-4-yl (1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethylcarbamate;~~

~~(2S)-4-cyclopropyl N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;~~

~~(2S)-4-cyclopentyl N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;~~

~~(1S)-1-cyclopropyl 2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl 4-methylpiperazine-1-carboxylate;~~

~~1-cyclopropyl 2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl-2-morpholin-4-ylethylcarbamate;~~

~~N-[(1S)-1-cyclopropyl 2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl]oxy-carbonyl]glycine;~~

~~(1S)-1-cyclopropyl 2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl 1-methylpiperidin-4-ylcarbamate;~~

~~(1S)-1-cyclopropyl 2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethylmethyl(1-methylpiperidin-4-yl)carbamate;~~

~~(1S)-1-cyclopropyl 2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl 4-dimethylamino)piperidine-1-carboxylate;~~

~~tert-butyl (2S)-4-(2-chloro-5-fluoropyrimidin-4-yl)-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxylate;~~

~~(2S)-4-(5-fluoro-2-methylpyrimidin-4-yl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;~~

~~(2S)-4-(2-chloro-5-fluoropyrimidin-4-yl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;~~

~~(2S)-4-(4-chloro-5-methylpyrimidin-2-yl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;~~

~~(2S)-4-(6-chloropyrimidin-4-yl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;~~

~~(2S)-4-(2-chloropyrimidin-4-yl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;~~

~~(2S)-N,N-dimethyl-4-(4-methylpyridin-3-yl)-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;~~

~~(2S)-N,N-dimethyl-2-phenyl-4-(1,3-thiazol-2-yl)-2,5-dihydro-1H-pyrrole-1-carboxamide;~~

~~(2S)-N,N-dimethyl-2-phenyl-4-(1,3-thiazol-4-yl)-2,5-dihydro-1H-pyrrole-1-carboxamide;~~

~~(2S)-N,N-dimethyl-2-phenyl-4-(1,2-thiazol-5-yl)-2,5-dihydro-1H-pyrrole-1-carboxamide;~~

~~4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N,N-dimethyl-2,5-dihydro-1H-pyrrole-1-carboxamide;~~

~~(2S)-4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N,N-dimethyl-2,5-dihydro-1H-pyrrole-1-carboxamide;~~

~~(2S)-4-(2,5-difluorophenyl)-N-(2-hydroxyethyl)-N-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;~~

~~N-{{[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}-N-methyl-beta-alanine};~~

~~methyl N-{{[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]carbonyl}-N-methyl-beta-alaninate};~~

~~4-{{[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]acetyl}morpholin-4-ium};~~

~~3-[(2S)-4-(2,5-difluorophenyl)-1-(2-hydroxy-2-methylpropanoyl)-2,5-dihydro-1H-pyrrol-2-yl]phenol;~~

~~4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N,N-dimethyl-2,5-dihydro-1H-pyrrole-1-sulfonamide;~~

~~3-[4-(2,5-difluorophenyl)-1-(methylsulfonyl)-2,5-dihydro-1H-pyrrol-2-yl]phenol;~~

~~3-[4-(2,5-difluorophenyl)-1-(morpholin-4-ylcarbonyl)-2,5-dihydro-1H-pyrrol-2-yl]phenol;~~

~~3-[4-(2,5-difluorophenyl)-1-(2,2-dimethylpropanoyl)-2,5-dihydro-1H-pyrrol-2-yl]phenol;~~

~~(2S)-4-(2,5-difluorophenyl)-1-[(methylsulfonyl)acetyl]-2-phenyl-2,5-dihydro-1H-pyrrole;~~

~~(2S)-4-(2,5-difluorophenyl)-2-phenyl-1-[(phenylsulfonyl)acetyl]-2,5-dihydro-1H-pyrrole;~~

~~3-[(2S)-1-[(2S)-2-cyclopropyl-2-hydroxyethanoyl]-4-(2,5-difluorophenyl)-2,5-dihydro-1H-pyrrol-2-yl]phenol;~~

~~3-[(2S)-4-(2,5-difluorophenyl)-1-[(2S)-2-hydroxy-3,3-dimethylbutanoyl]-2,5-dihydro-1H-pyrrol-2-yl]phenol;~~

~~(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethanol;~~

~~(2S)-4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-tetrahydrofuran-3-yl-2,5-dihydro-1H-pyrrole-1-carboxamide;~~

~~(2S)-4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-(2-methoxyethyl)-N-methyl-2,5-dihydro-1H-pyrrole-1-carboxamide;~~

~~(2S)-4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-tetrahydro-2H-pyran-4-yl-2,5-dihydro-1H-pyrrole-1-carboxamide;~~

~~3-[(2S)-4-(2,5-difluorophenyl)-1-(piperidin-1-ylcarbonyl)-2,5-dihydro-1H-pyrrol-2-yl]phenol;~~

4-(2,5-difluorophenyl)-N-[1-(2-fluoroethyl)piperidin-4-yl]-2-(3-hydroxyphenyl)-N-methyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-[[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]carbonyl](methyl)amino]piperidinium trifluoroacetate;

2-[[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]carbonyl](methyl)amino]ethyl 4-methylpiperazine-1-carboxylate;

3-{4-(2,5-difluorophenyl)-1-[(4-methylpiperazin-1-yl)carbonyl]-2,5-dihydro-1H-pyrrol-2-yl}phenol;

2-[[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]carbonyl](methyl)amino]ethyl morpholine-4-carboxylate;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[(1-methyl-1H-pyrazol-4-yl)methyl]-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-(isoxazol-5-ylmethyl)-N-methyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

2-[[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]carbonyl](methyl)amino]ethyl dimethylaminocarboxylate;

2-[[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]carbonyl](methyl)amino]ethyl piperidine-1-carboxylate;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[2-(2-oxopyrrolidin-1-yl)ethyl]-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[(5-oxo-4,5-dihydro-1H-1,2,4-triazol-3-yl)methyl]-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-(tetrahydro-2H-pyran-4-ylmethyl)-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-{{5-(methoxymethyl)-1H-pyrazol-3-yl)methyl}}-N-methyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-(1,3-thiazol-4-ylmethyl)-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[(4-methyl-1,2,5-oxadiazol-3-yl)methyl]-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-(1,3-thiazol-2-ylmethyl)-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-(isoxazol-3-ylmethyl)-N-methyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[2-(1H-1,2,4-triazol-1-yl)ethyl]-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[2-(1H-pyrazol-1-yl)ethyl]-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[(1-methyl-5-oxopyrrolidin-2-yl)methyl]-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-(1-isoxazol-3-ylethyl)-N-methyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-N-(1,3-dioxolan-2-ylmethyl)-2-(3-hydroxyphenyl)-N-methyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-N-(1,4-dioxan-2-ylmethyl)-2-(3-hydroxyphenyl)-N-methyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[(5-methyl-1,3,4-oxadiazol-2-yl)methyl]-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[2-(methylsulfonyl)ethyl]-2,5-dihydro-1H-pyrrole-1-carboxamide;

2-[{{4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl}carbonyl}(methyl)amino]ethanesulfonic acid;

~~2-hydroxyethyl (1S)-1-[[{(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl}carbonyl]-2,2-dimethylpropylcarbamate;~~

~~3-hydroxypropyl (1S)-1-[[{(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl}carbonyl]-2,2-dimethylpropylcarbamate;~~

~~2-hydroxyethyl {(1S)-1-isopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}carbamate;~~

~~2-hydroxyethyl {(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}carbamate;~~

~~4-hydroxybutyl (1S)-1-[[{(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl}carbonyl]-2,2-dimethylpropylcarbamate;~~

~~(2S)-4-(2,5-difluorophenyl)-1-[2-(methylsulfonyl)ethyl]-2-phenyl-2,5-dihydro-1H-pyrrole;~~



~~(2S)-4-(2,5-difluorophenyl)-1-[2-(ethylsulfonyl)ethyl]-2-phenyl-2,5-dihydro-1H-pyrrole;~~

~~1-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]pentan-3-one;~~

~~4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]butan-2-one;~~

~~4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3-methylbutan-2-one;~~

~~2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N,N-dimethylethanesulfonamide;~~

~~3-[(2S)-4-(2,5-difluorophenyl)-1-[2-(methylsulfonyl)ethyl]-2,5-dihydro-1H-pyrrol-2-yl]phenol;~~

~~methyl 3-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]propanoate;~~

~~(2S)-4-(2,5-difluorophenyl)-1-[2-(ethylsulfonyl)propyl]-2-phenyl-2,5-dihydro-1H-pyrrole;~~

~~3-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-methylpropanamide;~~

~~3-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N,N-dimethylpropanamide;~~

~~3-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N,N,2-trimethylpropanamide;~~

~~4-{3-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]propanoyl}morpholine;~~

~~1-{3-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]propanoyl}-4-(methylsulfonyl)piperazine;~~

~~1-{3-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]propanoyl}piperidin-4-ol;~~

~~methyl 3-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]propanoate;~~

~~2-[(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl]oxy-N-ethylacetamide;~~

~~4-[(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethoxy]acetyl}morpholine;~~

~~2-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethoxy)-N-(2-hydroxyethyl)acetamide;~~

~~1-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethoxy)acetyl)-4-methylpiperazine;~~

~~1-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethoxy)acetyl)piperazine;~~

~~2-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethoxy)-N-piperidin-4-ylacetamide;~~

~~1-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethoxy)acetyl)piperidin-4-amine;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-3-morpholin-4-yl-3-oxopropan-1-amine;~~

~~N<sup>3</sup>-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-N<sup>1</sup>,N<sup>1</sup>-dimethyl-β-alaninamide; and~~

~~((1S)-1-(((2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl)carbonyl)-2,2-dimethylpropyl)(3-morpholin-4-yl-3-oxopropyl)amine;~~

or a pharmaceutically acceptable salt or stereoisomer thereof.

12. (Currently amended) The compound according to Claim 11 which is selected from:

~~(-)-4-(2,5-difluorophenyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;~~

~~4-(5-chloro-2-fluorophenyl)-N,N-dimethyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;~~

~~4-(2,5-difluorophenyl)-1-(isopropylsulfonyl)-2-phenyl-2,5-dihydro-1H-pyrrole;~~

~~1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-methyl-1-oxopropan-2-ol;~~

~~1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-methyl-1-oxopropan-2-amine;~~

~~(1S)-1-[[[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-2,2-dimethylpropylamine; and~~

(2S)-4-(2,5-difluorophenyl)-N-methyl-2-phenyl-N-piperidin-4-yl-2,5-dihydro-1H-pyrrole-1-carboxamide;

or a pharmaceutically acceptable salt or stereoisomer thereof.

13. (Currently amended) The compound according to Claim 11 which is selected from:

~~(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethanamine;~~

(2S)-4-(2,5-difluorophenyl)-N-methyl-2-phenyl-N-piperidin-4-yl-2,5-dihydro-1H-pyrrole-1-carboxamide;

~~(1S)-1-[[[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-2,2-dimethylpropylamine;~~

~~2-(((1S)-1-tert-butyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}amino)-N-ethylacetamide;~~

(2S)-4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N,N-dimethyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(2,5-difluorophenyl)-N-methyl-N-(1-methylpiperidin-4-yl)-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide; and

~~(2S)-4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-1-L-valyl-2,5-dihydro-1H-pyrrole~~

or a pharmaceutically acceptable salt or stereoisomer thereof.

14. (Currently amended) A compound which is:

~~(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethanamine~~

~~(2S)-4-(2,5-difluorophenyl)-1-[(methylsulfonyl)acetyl]-2-phenyl-2,5-dihydro-1H-pyrrole~~

~~(2S)-4-(2,5-difluorophenyl)-2-phenyl-1-[(phenylsulfonyl)acetyl]-2,5-dihydro-1H-pyrrole~~

~~3-[(2S)-1-[(2S)-2-cyclopropyl-2-hydroxyethanoyl]-4-(2,5-difluorophenyl)-2,5-dihydro-1H-pyrrol-2-yl]phenol~~

~~3-[(2S)-4-(2,5-difluorophenyl)-1-[(2S)-2-hydroxy-3,3-dimethylbutanoyl]-2,5-dihydro-1H-pyrrol-2-yl]phenol~~

~~(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethanol~~

~~(2S)-4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-tetrahydrofuran-3-yl-2,5-dihydro-1H-pyrrole-1-carboxamide~~

~~(2S)-4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-(2-methoxyethyl)-N-methyl-2,5-dihydro-1H-pyrrole-1-carboxamide~~

~~(2S)-4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-tetrahydro-2H-pyran-4-yl-2,5-dihydro-1H-pyrrole-1-carboxamide~~

~~3-[(2S)-4-(2,5-difluorophenyl)-1-(piperidin-1-ylcarbonyl)-2,5-dihydro-1H-pyrrol-2-yl]phenol~~

~~4-(2,5-difluorophenyl)-N-[1-(2-fluoroethyl)piperidin-4-yl]-2-(3-hydroxyphenyl)-N-methyl-2,5-dihydro-1H-pyrrole-1-carboxamide~~

4-[[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]carbonyl](methylamino)piperidinium trifluoroacetate

2-[[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]carbonyl](methylamino)ethyl 4-methylpiperazine-1-carboxylate

3-{4-(2,5-difluorophenyl)-1-[(4-methylpiperazin-1-yl)carbonyl]-2,5-dihydro-1H-pyrrol-2-yl}phenol

2-[[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]carbonyl](methylamino)ethyl morpholine-4-carboxylate

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[(1-methyl-1H-pyrazol-4-yl)methyl]-2,5-dihydro-1H-pyrrole-1-carboxamide

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-(isoxazol-5-ylmethyl)-N-methyl-2,5-dihydro-1H-pyrrole-1-carboxamide

2-[[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]carbonyl](methylamino)ethyl dimethylaminocarboxylate

2-[[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]carbonyl](methylamino)ethyl piperidine-1-carboxylate

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[2-(2-oxopyrrolidin-1-yl)ethyl]-2,5-dihydro-1H-pyrrole-1-carboxamide

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[(5-oxo-4,5-dihydro-1H-1,2,4-triazol-3-yl)methyl]-2,5-dihydro-1H-pyrrole-1-carboxamide

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-(tetrahydro-2H-pyran-4-ylmethyl)-2,5-dihydro-1H-pyrrole-1-carboxamide

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N- {[5-(methoxymethyl)-1H-pyrazol-3-yl]methyl}-  
N-methyl-2,5-dihydro-1H-pyrrole-1-carboxamide

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-(1,3-thiazol-4-ylmethyl)-2,5-dihydro-  
1H-pyrrole-1-carboxamide

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[(4-methyl-1,2,5-oxadiazol-3-  
yl)methyl]-2,5-dihydro-1H-pyrrole-1-carboxamide

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-(1,3-thiazol-2-ylmethyl)-2,5-dihydro-  
1H-pyrrole-1-carboxamide

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-(isoxazol-3-ylmethyl)-N-methyl-2,5-dihydro-1H-  
pyrrole-1-carboxamide

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[2-(1H-1,2,4-triazol-1-yl)ethyl]-2,5-  
dihydro-1H-pyrrole-1-carboxamide

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[2-(1H-pyrazol-1-yl)ethyl]-2,5-dihydro-  
1H-pyrrole-1-carboxamide

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[(1-methyl-5-oxopyrrolidin-2-  
yl)methyl]-2,5-dihydro-1H-pyrrole-1-carboxamide

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-(1-isoxazol-3-ylethyl)-N-methyl-2,5-dihydro-1H-  
pyrrole-1-carboxamide

4-(2,5-difluorophenyl)-N-(1,3-dioxolan-2-ylmethyl)-2-(3-hydroxyphenyl)-N-methyl-2,5-dihydro-  
1H-pyrrole-1-carboxamide

4-(2,5-difluorophenyl)-N-(1,4-dioxan-2-ylmethyl)-2-(3-hydroxyphenyl)-N-methyl-2,5-dihydro-1H-pyrrole-1-carboxamide

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[(5-methyl-1,3,4-oxadiazol-2-yl)methyl]-2,5-dihydro-1H-pyrrole-1-carboxamide

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-[2-(methylsulfonyl)ethyl]-2,5-dihydro-1H-pyrrole-1-carboxamide

2-[[[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]carbonyl](methyl)amino]ethanesulfonic acid

2-hydroxyethyl (1S)-1-[[[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-2,2-dimethylpropyl]carbamate

3-hydroxypropyl (1S)-1-[[[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-2,2-dimethylpropyl]carbamate

2-hydroxyethyl {(1S)-1-isopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}carbamate

2-hydroxyethyl {(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}carbamate

4-hydroxybutyl (1S)-1-[[[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-2,2-dimethylpropyl]carbamate

(2S)-4-(2,5-difluorophenyl)-1-[2-(methylsulfonyl)ethyl]-2-phenyl-2,5-dihydro-1H-pyrrole

(2S)-4-(2,5-difluorophenyl)-1-[2-(ethylsulfonyl)ethyl]-2-phenyl-2,5-dihydro-1H-pyrrole

1-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]pentan-3-one

4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]butan-2-one

4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3-methylbutan-2-one

2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N,N-dimethylethanesulfonamide

3-[(2S)-4-(2,5-difluorophenyl)-1-[2-(methylsulfonyl)ethyl]-2,5-dihydro-1H-pyrrol-2-yl]phenol

~~methyl 3-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]propanoate~~

~~(2S)-4-(2,5-difluorophenyl)-1-[2-(ethylsulfonyl)propyl]-2-phenyl-2,5-dihydro-1H-pyrrole~~

~~3-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-methylpropanamide~~

~~3-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N,N-dimethylpropanamide~~

~~3-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N,N,2-trimethylpropanamide~~

~~4-{3-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]propanoyl}morpholine~~

~~1-{3-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]propanoyl}-4-(methylsulfonyl)piperazine~~

~~1-{3-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]propanoyl}piperidin-4-ol~~

~~methyl 3-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]propanoate~~

~~2-(((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}oxy)-N-ethylacetamide~~

~~4-(((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethoxy}acetyl)morpholine~~

~~2-(((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethoxy}-N-(2-hydroxyethyl)acetamide~~

~~1-(((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethoxy}acetyl)-4-methylpiperazine~~

~~1-(((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethoxy}acetyl)piperazine~~

~~2-(((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethoxy}-N-piperidin-4-ylacetamide~~

~~1-(((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethoxy}acetyl)piperidin-4-amine~~



~~*N*-{[(1*S*)-1-cyclopropyl-2-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-2-oxoethyl]-3-morpholin-4-yl-3-oxopropan-1-amine~~

~~*N*<sup>3</sup>-{[(1*S*)-1-cyclopropyl-2-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-2-oxoethyl]-*N*<sup>1</sup>,*N*<sup>1</sup>-dimethyl-β-alaninamide~~

~~[(1*S*)-1-[[[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]carbonyl]-2,2-dimethylpropyl](3-morpholin-4-yl-3-oxopropyl)amine~~

or a pharmaceutically acceptable salt thereof.

15. (Currently amended) The compound according to Claim 12 which is the TFA salt of a compound selected from:

~~2-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]sulfonyl]-*N,N*-dimethylethanamine;~~

~~1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-2-methyl-1-oxopropan-2-amine;~~

~~4-(5-chloro-2-fluorophenyl)-2-phenyl-1-(trifluoroacetyl)-2,5-dihydro-1*H*-pyrrole;~~

~~(1*S*)-1-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]carbonyl]-2-methylpropylamine;~~

~~(1*R*)-1-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]carbonyl]-2-methylpropylamine;~~

~~4-(2,5-difluorophenyl)-2-phenyl-1-*L*-prolyl-2,5-dihydro-1*H*-pyrrole;~~

~~4-(2,5-difluorophenyl)-2-phenyl-1-*D*-prolyl-2,5-dihydro-1*H*-pyrrole;~~

~~(4*R*)-4-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]carbonyl]-1,3-thiazolidine;~~

~~methyl (3S)-3-amino-4-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-4-oxobutanoate;~~

~~(4S)-4-amino-5-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-5-oxopentanamide;~~

~~(1S)-1-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-3-(methylthio)propylamine;~~

~~(1S)-1-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-3-(methylsulfonyl)propylamine;~~

~~(2S)-2-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]piperidine;~~

~~(1S)-1-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]pentylamine;~~

~~(1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxo-1-(thien-2-ylmethyl)ethylamine;~~

~~4-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-1,1-dioxidotetrahydro-2H-thiopyran-4-ylamine;~~

~~(2S)-1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-methyl-1-oxopropan-2-amine;~~

~~(1S)-1-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]propylamine;~~

~~(1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxo-1-phenylethanamine;~~

~~(1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxo-1-phenylethanamine;~~

~~(4S)-4-amino-5-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-5-oxopentanamide~~

~~3-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3-oxopropan-1-amine;~~

~~(1S,2S)-1-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-2-methylbutylamine;~~

~~(1S)-1-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]butylamine;~~

~~(1S)-1-cyclopropyl-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethanamine;~~

~~1-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]cyclopropanamine;~~

~~1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-oxopropan-2-amine;~~

~~(1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-methyl-2-oxoethylamine;~~

~~(1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxo-1-(pyridin-2-ylmethyl)ethylamine;~~

~~(1S)-1-cyclohexyl-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethanamine;~~

~~(1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-(4-iodobenzyl)-2-oxoethylamine;~~

~~(1S)-1-benzyl-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethylamine;~~

~~4-((2S)-2-amino-3-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3-oxopropyl)phenol;~~

~~(3S)-3-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-1,2,3,4-tetrahydroisoquinoline;~~

~~(1S)-1-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-3-phenylpropylamine;~~

~~(1S)-1-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-3-methylbutylamine;~~

~~(1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxo-1-(pyridin-3-ylmethyl)ethylamine;~~

~~1-[(2S)-azetidin-2-ylcarbonyl]-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrole;~~

~~(3S)-3-amino-4-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-4-oxobutanamide;~~

~~4-(2,5-difluorophenyl)-1-[(2-methylazetidin-2-yl)carbonyl]-2-phenyl-2,5-dihydro-1H-pyrrole;~~

~~(1S)-1-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-2,2-dimethylpropylamine;~~

~~methyl (4S)-4-amino-5-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-5-oxopentanoate;~~

~~4-(2,5-difluorophenyl)-2-phenyl-1-[(2S,3S)-2-phenylpyrrolidin-3-yl]carbonyl]-2,5-dihydro-1H-pyrrole;~~

~~4-(2,5-difluorophenyl)-2-phenyl-1-[(5-phenylpyrrolidin-3-yl)carbonyl]-2,5-dihydro-1H-pyrrole;~~

~~(2S)-2-amino-3-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3-oxopropan-1-ol;~~

~~(2R,3S)-3-amino-4-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-4-oxobutan-2-ol;~~

~~(1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-(methoxymethyl)-2-oxoethylamine;~~

~~4-(2,5-difluorophenyl)-2-phenyl-1-(pyrrolidin-3-ylcarbonyl)-2,5-dihydro-1H-pyrrole;~~

~~4-(2,5-difluorophenyl)-2-phenyl-1-[(3-phenylpyrrolidin-3-yl)acetyl]-2,5-dihydro-1H-pyrrole;~~

~~(1S)-1-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-3,3-difluoropropylamine;~~

~~(1S)-3-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3-oxo-1-phenylpropan-1-amine;~~

~~4-(2,5-difluorophenyl)-2-phenyl-1-[(4S)-4-phenyl-L-prolyl]-2,5-dihydro-1H-pyrrole;~~

~~1-{2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}cyclohexanamine;~~

~~2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethanamine;~~

~~4-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]piperidin-4-amine;~~

~~(1S,3R)-3-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]cyclopentanamine;~~

~~(1R,4S)-4-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]cyclopent-2-en-1-amine;~~

~~(1S,4R)-4-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]cyclopent-2-en-1-amine;~~

~~(1S)-1-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]but-3-ynylamine;~~

~~(1R)-3-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3-oxo-1-phenylpropan-1-amine;~~

~~3-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-2-phenylpiperidine;~~

~~(1S)-1-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]but-3-enylamine;~~

~~(2S)-3-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-(methylamino)-3-oxopropan-1-ol;~~

~~(3R,5S)-5-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]pyrrolidin-3-ol;~~

~~(1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxo-1-(1,3-thiazol-4-ylmethyl)ethylamine;~~

~~(1R)-1-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]but-3-enylamine;~~

~~(2S)-1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N,3-dimethyl-1-oxobutan-2-amine;~~

~~(2S)-1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N,4-dimethyl-1-oxopentan-2-amine;~~

~~(1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-[(1-methyl-1H-imidazol-4-yl)methyl]-2-oxoethylamine;~~

~~4-(2,5-difluorophenyl)-1-(N-6-formyl-L-lysyl)-2-phenyl-2,5-dihydro-1H-pyrrole;~~

~~(2S,3S)-1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N,3-dimethyl-1-oxopentan-2-amine;~~

~~(1S)-1-(cyclohexylmethyl)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethylamine;~~

~~(1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-(1H-indol-3-ylmethyl)-2-oxoethylamine;~~

~~(1S)-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-(isocyanomethyl)-2-oxoethylamine;~~

~~(1S)-1-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-3,3-dimethylbutylamine;~~

~~1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2,3-dimethyl-1-oxobutan-2-amine;~~

~~1-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]cyclohexanamine;~~

~~1-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]cyclopentanamine;~~

~~(1S)-3-(benzyloxy)-1-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]propylamine;~~

~~1-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2,3-dimethyl-1-oxobutan-2-amine;~~

~~1-[[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]cyclopent-3-en-1-amine;~~

~~(1S)-1-cyclopentyl-2-[4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethanamine;~~

~~4-(2,5-difluorophenyl)-1-(2-methylpropyl)-2-phenyl-2,5-dihydro-1H-pyrrole;~~

~~(1S)-2-[4-(5-chloro-2-fluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-1-cyclopropyl-2-oxoethanamine;~~

~~(1S,2S)-1-[[4-(5-chloro-2-fluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-2-methylbutylamine;~~

~~(1S)-1-[[[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]but-3-enylamine;~~

~~(1S)-1-[[[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]but-3-ynylamine;~~

~~(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethanamine;~~

~~1-cyclopropyl-3-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3-oxopropan-1-amine;~~

~~(1S,2S)-1-[[[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-2-methylbutylamine;~~

~~4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-1-L-valyl-2,5-dihydro-1H-pyrrole;~~

~~4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-1-(2-methylalanyl)-2,5-dihydro-1H-pyrrole;~~

~~3-[1-[(2S)-2-amino-2-cyclopropylethanoyl]-4-(5-chloro-2-fluorophenyl)-2,5-dihydro-1H-pyrrol-2-yl]phenol;~~



~~4-(5-chloro-2-fluorophenyl)-2-(3-hydroxyphenyl)-1-L-isoleucyl-2,5-dihydro-1H-pyrrole;~~

(2S)-4-(2,5-Difluorophenyl)-N-methyl-2-phenyl-N-(piperidin-4-ylmethyl)-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(5-Chloro-2-fluorophenyl)-N-methyl-2-phenyl-N-[(3R)-pyrrolidin-3-yl]-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(5-chloro-2-fluorophenyl)-N-methyl-2-phenyl-N-piperidin-4-yl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(5-chloro-2-fluorophenyl)-N-methyl-2-phenyl-N-[(3S)-pyrrolidin-3-yl]-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-4-(2,5-difluorophenyl)-N-methyl-2-phenyl-N-pyrrolidin-3-yl-2,5-dihydro-1H-pyrrole-1-carboxamide;

(2S)-N-(1-allylpiperidin-4-yl)-4-(2,5-difluorophenyl)-N-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-Difluorophenyl)-N-methyl-N-[(1-methylpiperidin-3-yl)methyl]-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-N-methyl-2-phenyl-N-(pyridin-3-ylmethyl)-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-N-methyl-N-[(1-methyl-1H-pyrazol-4-yl)methyl]-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-N-[2-(dimethylamino)ethyl]-N-methyl-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-N-methyl-2-phenyl-N-(2-pyridin-2-ylethyl)-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(5-chloro-2-fluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-piperidin-4-yl-2,5-dihydro-1H-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-N-methyl-N-piperidin-4-yl-2,5-dihydro-1H-pyrrole-1-carboxamide;

~~(2S)-1-[4-(2,5-difluorophenyl)-2-methyl-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-3-methyl-1-oxobutan-2-amine;~~

~~N-1-[(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl]-N-2,N-2-dimethylglycinamide;~~

~~N-1-[(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl]-N-2-methylglycinamide;~~

~~N-1-[(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}glycinamide;~~

~~N-1-[(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}-2-methylalaninamide;~~

~~N-[(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}-2-pyrrolidin-1-ylacetamide;~~

~~2-azetidin-1-yl-N-[(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl}acetamide;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-2-morpholin-4-ylacetamide;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-2-piperazin-1-ylacetamide;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-2-(4-methylpiperazin-1-yl)acetamide;~~

~~N-1-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-N-2-isopropylglycinamide;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)piperazine-1-carboxamide;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-N'-piperidin-4-ylurea;~~

~~4-amino-N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)piperidine-1-carboxamide;~~

~~N-(2-aminoethyl)-N'-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)urea;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-N'-(3-morpholin-4-ylpropyl)urea;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-N'-[2-(dimethylamino)ethyl]urea;~~

~~2-azetidin-1-yl-N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)ethanesulfonamide~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-2-(isopropylamino)ethanesulfonamide;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-2-pyrrolidin-1-ylethanesulfonamide;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-2-morpholin-4-ylethanesulfonamide;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-2-piperazin-1-ylethanesulfonamide;~~

~~N-((1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl)-2-(4-methylpiperazin-1-yl)ethanesulfonamide;~~

N-(tert-butyl)-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]acetamide;

2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-isopropylacetamide;

(2S)-1-(2-azetidin-1-yl-2-oxoethyl)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrole;

(2S)-4-(2,5-difluorophenyl)-1-(2-oxo-2-pyrrolidin-1-ylethyl)-2-phenyl-2,5-dihydro-1H-pyrrole;

4-{[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]acetyl}morpholine;

1-{[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]acetyl}piperazine;

1-{[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]acetyl}-4-methylpiperazine;

2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-isopropylbutanamide;

4-{2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]butanoyl}morpholine;

2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-ethylacetamide;

N-cyclobutyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]acetamide;

2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-ethylpropanamide;

N-cyclobutyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]propanamide;

2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-methylpropanamide;

2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-isopropylpropanamide;

N-(tert-butyl)-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]propanamide;

4-{2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]propanoyl}morpholine;

~~(3S)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-ethyl-2,2-dimethyl-4-oxobutanamide;~~

~~(3S)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2,2-dimethyl-4-oxo-N-piperidin-4-ylbutanamide;~~

~~(3S)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2,2-dimethyl-4-oxobutanoic acid;~~

~~(3S)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N,N,2,2-tetramethyl-4-oxobutanamide;~~

~~(1S)-1-[[[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl]-2,2-dimethyl-3-oxo-3-piperazin-1-yl]propylamine;~~

~~(3S)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N-isopropyl-2,2-dimethyl-4-oxobutanamide;~~

~~(3S)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N,2,2-trimethyl-4-oxobutanamide;~~

~~(3R)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-N,N,2,2-tetramethyl-4-oxobutanamide;~~

~~(3R)-3-amino-4-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2,2-dimethyl-4-oxobutanoic acid;~~

~~(1R)-1-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]carbonyl-2,2-dimethyl-3-oxo-3-piperazin-1-ylpropylamine~~

~~(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl-4-methylpiperazine-1-carboxylate;~~

~~(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl-1-methylpiperidin-4-ylcarbamate;~~

~~(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethylmethyl(1-methylpiperidin-4-yl)carbamate;~~

~~(1S)-1-cyclopropyl-2-[(2S)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1H-pyrrol-1-yl]-2-oxoethyl-4-dimethylamino)piperidine-1-carboxylate;~~

~~(2S)-N,N-dimethyl-4-(4-methylpyridin-3-yl)-2-phenyl-2,5-dihydro-1H-pyrrole-1-carboxamide;~~

~~4-[[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1H-pyrrol-1-yl]acetyl]morpholin-4-ium;~~

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-*N*-methyl-*N*-piperidin-4-yl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

2-[[4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-2,5-dihydro-1*H*-pyrrol-1-yl]carbonyl](methyl)amino]ethyl-4-methylpiperazine-1-carboxylate;

3-{4-(2,5-difluorophenyl)-1-[(4-methylpiperazin-1-yl)carbonyl]-2,5-dihydro-1*H*-pyrrol-2-yl}phenol;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-*N*-methyl-*N*-[(1-methyl-1*H*-pyrazol-4-yl)methyl]-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

4-(2,5-difluoro-phenyl)-2,5-dihydro-2-(3-hydroxyphenyl)-*N*-methyl-*N*--[2-(4-methyl-1-piperazinyl)-2-oxoethyl]-*H*-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-*N*-methyl-*N*-[(5-oxo-4,5-dihydro-1*H*-1,2,4-triazol-3-yl)methyl]-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-*N*-{[5-(methoxymethyl)-1*H*-pyrazol-3-yl)methyl}-*N*-methyl-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-*N*-methyl-*N*-(1,3-thiazol-4-ylmethyl)-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-*N*-methyl-*N*-(1,3-thiazol-2-ylmethyl)-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-*N*-methyl-*N*-[2-(1*H*-1,2,4-triazol-1-yl)ethyl]-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

4-(2,5-difluorophenyl)-2-(3-hydroxyphenyl)-*N*-methyl-*N*-[2-(1*H*-pyrazol-1-yl)ethyl]-2,5-dihydro-1*H*-pyrrole-1-carboxamide;

~~4-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]butan-2-one;~~

~~4-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-3-methylbutan-2-one;~~

3-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-*N*-methylpropanamide;

3-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-*N,N*-dimethylpropanamide;

3-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]-*N,N*,2-trimethylpropanamide;

4-{3-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]propanoyl}morpholine;

1-{3-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]propanoyl}-4-(methylsulfonyl)piperazine;

1-{3-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]propanoyl}piperidin-4-ol; and

methyl 3-[(2*S*)-4-(2,5-difluorophenyl)-2-phenyl-2,5-dihydro-1*H*-pyrrol-1-yl]propanoate.

16. (original) A pharmaceutical composition that is comprised of a compound in accordance with Claim 1 and a pharmaceutically acceptable carrier.

17. – 42. (Cancelled)